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# ROY F. WESTON, INC.

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REMEDIAL INVESTIGATION REPORT EVERGREEN MANOR SITE ROSCOE, ILLINOIS





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28 March 2001

Mr. Mike Ribordy Remedial Project Manager U.S. Environmental Protection Agency 77 West Jackson Boulevard Chicago, Illinois 60604

U.S. EPA Contract No.: 68-W7-0026 Work Assignment No.: 036-RICO-05MZ Document Control No.: RFW036-2E-AHVH

Re: Remedial Investigation Report

Evergreen Manor Site

Roscoe, Illinois

Dear Mr. Ribordy:

Roy F. Weston. (WESTON®) is pleased to submit three copies of the Final Remedial Investigation (RI) Report for the Evergreen Manor site. If you have any questions, please contact us at (847) 918-4000.

Very truly yours,

ROY F. WESTON, INC.

Kurt T. Fischer, P.G.

Site Manager

KTF:sk

**Enclosures** 

## REMEDIAL INVESTIGATION REPORT EVERGREEN MANOR SITE ROSCOE, ILLINOIS

**VOLUME I** 

March 2001

Prepared for
U.S. Environmental Protection Agency
77 West Jackson Boulevard
Chicago, Illinois 60604

This document was prepared in accordance with U.S. EPA Contract No. 68-W7-0026, WESTON Region V Response Action Contract (RAC) and contains confidential business information.

#### REMEDIAL INVESTIGATION REPORT EVERGREEN MANOR SITE ROSCOE, ILLINOIS

U.S. EPA CONTRACT NO. 68-W7-0026

Work Assignment No. 036-RICO-05MZ Document Control No. RFW036-2A-AHVH

March 2001

Prepared

and

Kurt Fischer

Approved By:

Site Manager

Approved By

James M. Burton, P.É.

Program Manager

#### **TABLE OF CONTENTS**

Sectio	<u>n</u> <u>Title</u>	<u>Page</u>
<u>volu</u>	UME I	
1	INTRODUCTION	1-1
	1.1 Objectives and Scope of Remedial Investigation	1-1
	1.2 Report Organization	1-2
2	SITE BACKGROUND	2-1
	2.1 Site Description	2-1
	2.2 Site History and Previous Reports	2-2
3	ENVIRONMENTAL SETTING	3-1
	3.1 Demography and Land Use	3-1
	3.2 Climate	3-1
	3.3 Ecology	3-2
	3.4 Regional Topography	3-3
	3.5 Regional Geology	3-3
	3.6 Soils	3-4
	3.7 Regional Hydrogeology	3-5
	3.8 Regional Surface Water Hydrology	3-7
	3.9 Groundwater Use in the Area	3-7
4	ENVIRONMENTAL INVESTIGATION PROCEDURES	4-1
	4.1 Fracture Trace Analysis	4-2
	4.2 Cone Penetrometer Testing (CPT)	4-2
	4.3 CPT Groundwater Sampling	4-4
	4.4 Monitoring Well Sampling	4-5
	4.5 Residential Well Sampling	4-6
	4.6 Sediment Sampling	4-7
	4.7 Surface Water Sampling	4-8
	4.8 Groundwater Elevation Measurements	4-8
5	ENVIRONMENTAL INVESTIGATION RESULTS	5-1
	5.1 Fracture Trace Analysis	5-1
	5.2 CPT Groundwater Sampling	5-2
	5.3 Monitoring Well Sampling	5-4
	5.4 Residential Well Sampling	5-6
	5.5 Sediment Sampling	5-8
	5.6 Surface Water Sampling	5-9
	5.7 Groundwater Elevation Measurements	5-9

CHO1\PUBLIC\WO\RAC\036\29672TOC.WPD

RFW036-2A-AHVH

## **TABLE OF CONTENTS (Continued)**

<u>Sectio</u>	<u>on</u>	<u>Title</u>	Page
6	GEOLOG	6-1	
	6.1 Si	ite Geology	6-1
	6.	1.1 CPT Interpretation	6-1
	6.2 Si	ite Hydrogeology	6-3
	6.3 Si	ite Hydrology	6-4
. 7	NATURI	7-1	
	7.1 D	7.1 Development of Screening Levels	
	7.	1.1 Groundwater	7-1
		1.2 Sediment	7-2
	7.	1.3 Surface Water	7-2
	7.2 Sc	ources of Contamination	7-3
		xtent of Contamination in Groundwater	7-3
	7.4 E	xtent of Contamination in Sediment	7-4
	7.5 Ex	xtent of Contamination in Surface Water	7-4
8	CONTAI	8-1	
	8.1 C	ontaminants of Potential Concern	8-1
	8.2 Fa	ate of Contaminants of Concern	8-2
	8.3 C	ontaminant Migration Pathway	8-3
9	HUMAN HEALTH RISK ASSESSMENT		9-1
	9.1 In	troduction	9-1
	9.	1.1 Objectives	9-1
	9.	1.2 Risk Assessment Approach	9-2
	9.	1.3 Risk Assessment Organization	9-2
	9.2 C	ontaminant Identification	9-3
	9.	2.1 Contaminant Characterization	9-3
	9.	2.2 Data Evaluation	9-5
	9.	2.3 Identification of Constituents of Potential Concern	9-6
	9.3 Ex	kposure Assessment	9-7
		3.1 Characterization of Exposure Setting	9-8
	9.	3.2 Identification of Exposure Pathways	9-8
	9.	3.3 Quantification of Exposure	9-11
		oxicity Assessment	9-16
		isk Characterization	9-17
		5.1 Quantitative Evaluation for Chemical COPCs	9-17
		5.2 Residential Scenario	9-19
		5.3 Commercial/Industrial Worker Scenario	9-20
	9.	5.4 Uncertainty Analysis	9-21

ii

## TABLE OF CONTENTS (Continued)

Section			<u>Title</u>	
10	ECO	LOGICA	L RISK ASSESSMENT	10-1
	10.1	Screeni	ing-Level Problem Formulation	10-2
		10.1.1	Environmental Setting	10-2
		10.1.2	Extent of Contamination	10-4
		10.1.3	Contaminant Fate and Transport	10-4
		10.1.4	Potential Ecological Receptors	10-4
		10.1.5	Complete Exposure Pathways	10-5
		10.1.6	General Assessment Endpoints	10-5
		10.1.7	Conceptual Model	10-6
	10.2	Screeni	ing-Level Ecological Effects Evaluation	10-6
		10.2.1	Constituents of Potential Concern	10-6
		10.2.2	Toxicity Reference Values	10-6
10.3		Screeni	ing-Level Exposure Estimate	10-7
	10.4 Screening-Level Risk Characterization		10-8	
		10.4.1	Summary of Uncertainty	10-8
		10.4.2	Risk Description	10-9
11	CON	CLUSIO	NS AND RECOMMENDATIONS	11-1
	11.1	11.1 Conclusions		11-1
		11.1.1	Site Geology and Hydrogeology	11-1
		11.1.2	Nature and Extent of Contamination	11-2
		11.1.3	Contaminant Fate and Transport	11-4
		11.1.4	Risk Assessment	11-5
		11.1.5	Contaminant Concentrations	11-6
		11.1.6	Non-Time Critical Removal Action	11-6
	11.2	Recomm	mendations	11-7
12	REFE	RENCES	S	12-1

#### **LIST OF FIGURES**

<u>Figure</u>	<u>Title</u>	
2-1	Site Layout	2-7
4-1	Groundwater Sampling Locations (CPT and Monitoring Wells)	4-9
4-2	Residential Well Location Map	4-10
4-3	Surface Water and Sediment Sampling Location Map	4-11
6-1	Cross Section Transect Map	6-6
6-2	Geologic Cross Section A - A'	6-7
6-3	Geologic Cross Section B - B'	6-8
6-4	Geologic Cross Section C - C'	6-9
6-5	Groundwater Piezometric Surface Contour Map	6-10
7-1	Extent of Positive Detections in Groundwater	7-5
10-1	Wetlands Near Evergreen Manor Site	10-11

### LIST OF TABLES

<u>Table</u>	<u>Title</u>	
5-1	Groundwater Analytical Results - CPT - VOCs	5-10
5-2	Groundwater Analytical Results - Monitoring Wells - VOCs	5-22
5-3	Groundwater Analytical Results - Monitoring Wells -	
	Water Quality Parameters	5-25
5-4	Groundwater Analytical Results - Residential Wells - VOCs	5-26
5-5	Groundwater Analytical Results - Residential Wells -	
	Water Quality Parameters	5-30
5-6	Sediment Analytical Results - VOCs	5-31
5-7	Sediment Analytical Results - Total Organic Carbon	5-33
5-8	Surface Water Analytical Results - VOCs	5-34
6-1	Groundwater Elevations from Monitoring Wells	6-11
7-1	Screening Levels for Detected Constituents	7-6
7-2	Summary of Detected Constituents in Groundwater - CPT-VOCs	7-7
7-3	Summary of Detected Constituents in Groundwater -	
	Monitoring Wells-VOCs	7-15
7-4	Summary of Detected Constituents in Groundwater -	
	Residential Wells - VOCs	7-17
7-5	Summary of Detected Constituents in Sediments - VOCs	7-20
8-1	Physical and Chemical Properties of Organic Constituents of Potential	
	Concern	8-11
8-2	Migration Characteristics of Constituents of Potential Concern	8-12
8-3	Comparison of Previous Groundwater Data with RI Data	8-13
8-4	Comparison of Select Detected Constituents in Groundwater - Residential	
	Wells - VOCs	8-16
8-5	Comparison of Select Detected Constituents in Groundwater - Monitoring	
	Wells - VOCs	8-17
8-6	BIOSCREEN Model Inputs and Results	8-20
8-7	BIOSCREEN Calibration Data	8-21
9-1	Permeability Coefficients for Organic COPCs	9-25
9-2	Total Carcinogenic Risk Associated with Chemical COPC Exposure	9-26
9-3	Total Hazard Index Associated with Chemical COPC Exposure	9-27
9-4	Summary of Uncertainty Analysis	9-28
10-1	Preliminary Ecological Conceptual Site Model	10-12
10-2	Comparison of Groundwater Concentrations to Surface Water	
	Screening Values	10-13

CHO1\PUBLIC\WO\RAC\036\29672TOC.WPD

#### APPENDICES

Appendix A - Supporting Material for Human Health and Ecological Risk Assessments

Appendix B - Fracture Trace Analysis, Resolution Resources, Inc.

#### **VOLUME II**

Appendix C - Water Sample Collection Forms

Appendix D - Analytical Data

Appendix E - Report for Cone Penetrometer Testing Services, Fugro Geosciences, Inc.

Appendix F - BIOSCREEN Model Results

Section: 1 Revision: 0

Date: March 2001 Page: 1 of 3

**SECTION 1** 

INTRODUCTION

This Remedial Investigation (RI) report presents the results from a field investigation conducted at

the Evergreen Manor site, in Roscoe, Illinois (hereafter referred to as the Evergreen Manor site). The

RI field investigation was conducted by Roy F. Weston, Inc. (WESTON<sub>®</sub>) for the United States

Environmental Protection Agency (U.S. EPA) from May 2000 to June 2000. This RI Report was

prepared in accordance with the U.S. EPA Guidance for Conducting Remedial Investigations and

Feasibility Studies under CERCLA, Interim Final (U.S. EPA, 1988).

1.1 OBJECTIVE AND SCOPE OF REMEDIAL INVESTIGATION

In accordance with the U.S. EPA guidance document (U.S. EPA, 1988), the objective of the RI is

to gather site information sufficient to support a Feasibility Study (FS) and to make an informal risk

management decision regarding an appropriate remedy. The specific objectives of the RI are the

following:

• Extent of Contamination: To evaluate the areal and vertical extent of

contamination.

• Contaminant Fate and Transport: To determine the rate of migration and the fate

of contaminants through various migration pathways.

Ecological Assessment: To characterize and estimate the potential for adverse

ecological effects associated with the contamination at the site.

Baseline Risk Assessment: To evaluate the potential human health and

environmental impacts associated with the site under a no-action alternative (i.e., in

the absence of remedial action).

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RFW036-2A-AHVH

Revision: 0

Date: March 2001 Page: 2 of 3

The RI was conducted in accordance with the Evergreen Manor, Quality Assurance Project Plan

(QAPP) and Field Sampling Plan (FSP) (WESTON, 2000a). The main RI investigation tasks

consisted of a fracture trace analysis, cone penetrometer testing (CPT), groundwater sampling,

sediment sampling, surface water sampling, and monitoring well sampling.

1.2 REPORT ORGANIZATION

This RI report is divided into the following sections:

Executive Summary: Presents an overall summary of the RI Report.

Section 1, Introduction: Provides a brief overview of the objective and scope of the RI.

<u>Section 2, Site Background</u>: Provides the site description, site history, and a summary of results of previous site investigations.

Section 3, Environmental Setting: Describes the surrounding land use and population, climate, regional soils, surface water features, area drainage and topography, regional geology, regional hydrogeology, and groundwater use in the area.

<u>Section 4, Environmental Investigation Procedures</u>: Describes the procedures of the RI field investigation.

<u>Section 5, Environmental Investigation Results</u>: Describes the results of the RI field investigation.

<u>Section 6, Results of Geologic/Hydrogeologic Investigations</u>: Provides descriptions of the site geology and hydrogeology based on the results of the RI field investigation.

Section 7. Nature and Extent of Contamination: Presents the nature and extent of contamination by each medium.

<u>Section 8, Contaminant Fate and Transport</u>: Describes the contaminant migration pathways, environmental fate of contaminants, and rate of contaminant migration.

<u>Section 9, Human Health Risk Assessment:</u> Provides an evaluation of the potential threat to human health and the environment in the absence of any remedial actions.

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RFW036-2A-AHVH

Section: 1 Revision: 0 Date: March 2001 Page: 3 of 3

<u>Section 10, Ecological Assessment Summary</u>: Provides an evaluation of the potential impacts to the ecological community from site-related contaminants.

<u>Section 11, Conclusions and Recommendations</u>: Presents the conclusions of the RI, identifies data gaps, and makes recommendations for additional work.

Section 12, References: Lists all reference sources used in the RI report.

The tables and figures are referenced by section numbers and are presented at the end of each section.

Section: 2 Revision: 1

Date: 28 March 2001

Page: 1 of 7

**SECTION 2** 

SITE BACKGROUND

2.1 SITE DESCRIPTION

The Evergreen Manor site is located approximately 1.5 miles northwest of Roscoe, Winnebago

County, Illinois. The site includes four residential subdivisions and has been defined by the extent

of groundwater contamination. A site layout is provided in Figure 2-1. The site extends over

Sections 16, 20, 21, 29, and 32 in Township 46 North, Range 2 East, and is found on the South

Beloit, Illinois/Wisconsin Quadrangle. The coordinates of the site are latitude 42° 26' 32.0".

longitude 89° 01' 36.0".

The site area was used as farmland prior to development. The Hononegah Heights subdivision was

developed between 1940 and 1964; the Tresemer subdivision was developed between 1972 and

1974; the Olde Farm subdivision was developed between 1976 and 1979; and the Evergreen Manor

subdivision was developed between 1986 and 1988. With the exception of the Evergreen Manor

subdivision, most of the development occurred in the late 1970's and the early 1980's.

The Evergreen Manor site is bounded to the south by the Rock River. The Hononegah Forest

preserve is located to the west of the site, some agricultural fields are located to the east of the site,

and agricultural land is located north of the site (IEPA, 1992).

Roscoe Rock and Sand, Inc., a gravel pit and concrete mixing facility, is located approximately 0.5

miles to the northeast of the site. Roscoe Rock and Sand, Inc. purchased the former Kelley Sand and

Gravel property, and is located on the north and south sides of McCurry Road, west of Route 251.

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RFW036-2A-AHVH

Section: 2 Revision: 1

Date: 28 March 2001

Page: 2 of 7

An industrial park is located approximately 2 miles to the northeast of the site and is located north

of Rockton Road, and east of Route 251. The industrial park contains the following businesses:

Inlander-Steindler Paper Company, Regal-Beloit Corporation, McGuire Brothers Auto Body and

Sand Blasting, Makerite Manufacturing Company, Midwest-Precision Grinding, Rockford Steam

Boiler Works, Oscar's Auto and Battery Clinic, Dayles Welding, Armour Specialty, Inc. (industrial

painting), RD Systems, Electro Cam Corporation, Area Elevator, DGM, Preston 151 (trucking firm),

and Indicon Midwest (IEPA, 1992).

Several industries are located on the south side of Rockton Road, east of Illinois Route 251. These

include Ecolab and Taylor Design, Inc. Further south along and east of Illinois Route 251 are five

other facilities: State Line Foundries, Waste Management Transfer Station, Kenny's Cars, Trucks

and Equipment, Stateline Printing Company, and Stateline Storage. Warner Brake and Clutch is

located south of McCurry Road on the east side of Route 251 (IEPA, 1992).

2.2 SITE HISTORY AND PREVIOUS REPORTS

The Evergreen Manor site was first realized in November 1990 when a lending institution required

a homeowner to sample and analyze the drinking water. Results of the analyses indicated elevated

concentrations of volatile organic compounds (VOCs). The Illinois Department of Public Health

(IDPH) undertook further sampling in the area and identified a plume of contamination located

beneath the Hononegah Heights and Evergreen Manor subdivisions (IEPA, 1992).

The results of the IDPH sampling indicated that maximum concentration limits (MCLs) were

exceeded for trichloroethene (TCE; MCL of 5 ppb), and 1,1-dichloroethene (1,1-DCE; MCL of 7

ppb) in one or more wells. Other VOCs identified included cis-1,2-dichloroethene (c-1,2-DCE),

1,1-dichloroethane (1,1-DCA), 1,1,1-trichloroethane (1,1,1-TCA), tetrachloroethane (PCE), and

1,1,2-trichloroethane (1,1,2-TCA). Based on these results, the IDPH concluded that at least 130

CH01\PUBLIC\WO\RAC\036\29672S-2.WPD

RFW036-2A-AHVH

Section: 2 Revision: 1

Date: 28 March 2001

Page: 3 of 7

residences in the Hononegah, Olde Farm, Evergreen Manor, and possibly the Tresemer subdivisions

could be contaminated with VOCs (IEPA, 1992).

The Evergreen Manor site was added to the Comprehensive Environmental Response, Compensation

and Liability Information System (CERCLIS) on 3 August 1991, based on information received by

the Illinois Environmental Protection Agency (IEPA) from the IDPH and the U.S. Environmental

Protection Agency (U.S. EPA). The first Comprehensive Environmental Response and Liability Act

(CERCLA) evaluation was a Preliminary Assessment (PA) of the Evergreen Manor site in January

1992.

CERCLA Screening Site Inspection Report, 1992

A Site Screening Inspection (SSI) was performed in June and August 1992 to gather information for

potential Hazard Ranking. A total of 39 soil gas samples and 4 groundwater samples were collected

and analyzed for 1,1,1-TCA, TCE, and 1,1-DCE. Soil gas samples collected along McCurry Road,

east of IL Route 251, and along the frontage road east of IL Route 251 indicated the presence of the

three VOCs. The compounds were not detected on the north side of the Ecolab facility, nor along

the north side of Rockton Road.

The results of the groundwater analyses indicated that the VOCs were not detected in the samples

collected on the north side of the Ecolab facility, or along the frontage road. The three VOCs were

detected in a well north of the Waste Management facility.

The SSI Report assigned a high priority to the Evergreen Manor site based on the results of the SSI

and the groundwater samples collected from residential wells by IDPH between 1990 and December

1991.

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RFW036-2A-AHVH

Section: 2 Revision: 1

Date: 28 March 2001

Page: 4 of 7

CERCLA Expanded Site Inspection Report, 1999

The Expanded Site Inspection (ESI) was conducted in November 1993 and consisted of the

collection of 49 groundwater samples from residential wells. The residential wells sampled are

located in the Hononegah Heights, Olde Farm, and Evergreen Manor subdivisions. The purpose of

the ESI was to collect information in support of the Hazard Ranking System package preparation.

A total of 49 groundwater samples were collected from 45 residences in the three subdivisions, four

of which were duplicate samples. The samples were analyzed for VOCs. Results indicated that in

all but one of the samples, and excluding two background samples, at least one VOC was detected.

Acetone, 1,1-DCE, 1,1-DCA, 1,2-DCE (total), and PCE were detected at concentrations less than

the laboratory detection limits. 1,1,1-TCA and TCE were detected at concentrations of less than 10

ppb to 37 and 40 ppb, respectively. Of the 45 wells sampled, 36 were found to have 1,1,1-TCA

concentrations significantly above background, and 40 were found to have TCE concentrations

significantly above background. All 40 TCE detections were at concentrations greater than the

MCL.

Additionally, this report indicated that results of residential well sampling by IEPA in December

1993 and January 1994 found more than 60 residential wells with concentrations of 1,1,1-TCA

and/or TCE above either the MCLs or the Cancer Risk.

Hazard Ranking System Documentation Record, 29 May 1997

Based on the information and data gathered from the PA, the SSI, the ESI, and other sampling by

IEPA and/or IDPH, a Hazard Ranking System (HRS) score was prepared. The Evergreen Manor

Ground Water Contamination Plume, ILD 984836734, received a score of 100 points, for the

groundwater pathway. The air, soil, and surface water pathways were not evaluated. The final HRS

site score was 50 points.

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RFW036-2A-AHVH

Section: 2 Revision: 1

Date: 28 March 2001

Page: 5 of 7

Engineering Evaluation/Cost Analysis Report, October 1998

The Engineering Evaluation/Cost Analysis Report (EE/CA) was written with the objective of

evaluating removal action objectives and removal action alternatives. Three viable alternatives were

identified which would abate the threat posed by drinking the groundwater. These included a

point-of-entry carbon filter treatment option, a point-of-use carbon filter treatment option, and an

option to connect the affected residences to a public water supply system.

The EE/CA also summarized work that has occurred at the Evergreen Manor site outside of the

scope of the PA, SSI, and ESI:

• IEPA and IDPH sampled 267 drinking water wells, mostly in the four subdivisions,

between December 1990 and March 1994. Results indicated that 108 wells exceeded

MCLs and 203 were impacted.

• IEPA installed 24 monitoring wells between December 1993 and February 1995.

Sample results from March 1994 indicated that 2 out of 20 wells exceeded MCLs

for TCE and PCE. Sample results from February 1995 indicated that 3 out of 24

wells exceeded MCLs for TCE and four exceeded for PCE.

Sample results from 12 wells sampled by U.S. EPA on 22 May 1998 indicated that

six wells exceeded the MCLs for TCE and three wells exceeded the MCLs for PCE.

Action Memorandum, 2 March 1999

This Action Memorandum served as a request for a non-time critical CERCLA removal action and

consistency exemption to the \$2 million and 12 month statutory limit at the Evergreen Manor site.

The U.S. EPA recommended the extension of the public water supply system in order to provide the

affected residences with safe drinking water. This decision was based on the permanence of the

solution and the public response to the EE/CA during the public comment period, which opened on

10 November 1998.

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RFW036-2A-AHVH

Section: 2 Revision: 1

Date: 28 March 2001

Page: 6 of 7

The Evergreen Manor site contamination was estimated to affect 700 people in approximately 250

residences. A preliminary risk assessment indicated that the continued usage of residential wells

would pose a threat to public health and the environment. Since the concentrations of TCE and PCE

exceed MCLs, this condition represented an imminent and substantial endangerment to local

residents.

The source area was identified in the Action Memorandum as the area at the intersection of Rockton

Road and Route 251.

Administrative Order on Consent, 14 May 1999

U.S. EPA entered into an administrative order on consent (AOC) with three PRPs concerning the

Evergreen Manor Groundwater Contamination Site. The AOC requires the PRPs to pay a total of

\$2,100,850 to partially fund the removal action to be performed by U.S. EPA. The removal action

will consist of construction of a water main extension to bring potable water from the North Park

Public Water District to the individual residences threatened by contaminated water.

Work related to the extension of the public water supply system and hookup of the effected

residences commenced in 1999, and was completed on September 29, 2000.

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BASEMAP ADAPTED FROM FIGURE 1 OF THE CONTAMINANT SOURCE EVALUATION — EVERGREEN MANOR SITE, DATED JANUARY 1997, BY CONESTOGA—ROVERS & ASSOCIATES.

750 E. Bunker Ct. SITE LAYOUT
Suite 500
Vernon Hills, Illinois
60061
Roscoe, Illinois

Evergreen Manor

Remedial Investigation Report

Section: 3 Revision: 1

Date: 28 March 2001

Page: 1 of 7

**SECTION 3** 

**ENVIRONMENTAL SETTING** 

3.1 DEMOGRAPHY AND LAND USE

The Evergreen Manor site in Roscoe, Winnebago County, Illinois includes four residential

subdivisions and has been defined by the extent of groundwater contamination. A site layout is

provided in Figure 2-1. According to the 1990 U.S. Census, the Evergreen Manor site is in Census

Tract 003902, Block Group 2, and has a total population of 3,632, 3% of whom are minorities. The

Evergreen Manor contamination has the potential for affecting 243 homes and an estimated

population of approximately 700 persons (IDPH, 1999).

Land use in and around the site is residential, agricultural, commercial and industrial. The land from

Hononegah Road to the Rock River is residential. Directly north of Hononegah Road, is commercial

property with various stores in a strip mall. Between the strip mall and Rockton Road, on the west

side of Illinois Route 251, most of the land is agricultural and is actively used during the growing

season. This area also includes an area of heavy industrial land use: a sand and gravel quarry and

cement mixing facility. To the east of Illinois Route 251, from Hononegah Road to Rockton Road,

land use is mixed between commercial, light industrial and residential. This area includes the Ecolab

facility, the Waste Management Transfer Station, Kenny's Cars, Trucks, and Equipment, as well as

other companies. In the northeast quadrant of Illinois Route 251 and Rockton Road most of the land

is light industrial and is occupied by the industrial park. Section 2.1 lists all of the industries and

businesses in the site area.

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RFW036-2A-AHVH

Section: 3 Revision: 0

Date: 31 October 2000

Page: 2 of 7

3.2 CLIMATE

Winnebago County has a continental climate typical of northern Illinois. This area has hot summers

and cold winters, with July being the hottest month and January being the coldest. The average

temperature in winter is 23°F and in summer is 71°F. The lowest recorded temperature was -22 °F

recorded on 21 January, 1970. The highest recorded temperature was 103 °F recorded on 27 July,

1955. Annual precipitation averages 38 inches and annual snowfall averages 33 inches. Sixty-six

percent of the rainfall occurs between the months of April through September.

3.3 ECOLOGY

The site is located in the Central Lowland geomorphic province, in the eastern broadleaf forest

province of the Hot Continental Division in the Humid Temperate Domain (USDA Forest Service,

Ecological Subregions of the United States, http://www.fs.fed.us/land/pubs/ecoregions).

The Rock River receives drainage from three major streams - the Pecatonica River, the Kishwaukee

River, and the Green River. It is 163 miles long in Illinois, and drains 2,272,000 acres in Illinois.

Of the total river miles in this basin, 69 stream miles have "good" overall resource quality and 97.9

miles have "fair" quality. The Rock River enters the Mississippi River at Rock Island (IDNR,

http://dnr.state.il.us/lands/education/valerie/end/page6.htm). At Rockton, the mean daily discharge

ranges from 2839 cubic feet per second (cfs) in September to 7375 cfs in April, with an annual mean

of 4178 cfs (USGS, CD-ROM, Current Year Discharge, http://www.il.water.usgs.gov/cd04-

99/dis tbl/05437500.htm). Dry Creek, a tributary of the Rock River, enters the river northwest of

the Tresemer Subdivision. West of Dry Creek, the river is classified by the NWI as a riverine

wetland and east of the creek, the river is classified as a lacustrine system. Forested wetlands border

both the river and the creek west of the site and the river south of the site. There are small areas of

emergent wetlands within the Evergreen Manor subdivision.

CH01\PUBLIC\WO\RAC\036\29672S-3.WPD

RFW036-2A-AHVH

Section: 3 Revision: 0

Date: 31 October 2000

Page: 3 of 7

The U.S. Fish and Wildlife Service (USFWS) was contacted to obtain information on threatened and

endangered species within the Evergreen Manor project area. Species that may be present in the area

include the endangered Indiana bat (Myotis sodalis), the threatened prairie bush clover (Lespedeza

leptostachya), the threatened bald eagle (Haliaeetus leucocephalus). Further information on these

species and their critical habitat is provided in Appendix A.

3.4 REGIONAL TOPOGRAPHY

The topography in Winnebago County has been created in large part by features developed during

the advance and retreat of glaciers. This includes till plains that contain kames, drumlins, and eskers

(USDA-NRCS, 1980). The Evergreen Manor site is located on a broad, flat terrace, which gently

slopes toward the Rock River. Locally, relief is no greater than about 70 feet from the highest area

near Rockton Road and IL Route 251, down to the Rock River elevation of approximately 700 feet

above Mean Sea Level (MSL).

3.5 REGIONAL GEOLOGY

The geology in the vicinity of the Evergreen Manor site has been most heavily influenced by fluvial

and glacial processes. The preglacial Rock River incised a deep bedrock valley that was

subsequently buried during glaciation. As the glacier retreated, vast deposits of sand and gravel with

lesser amounts of silt and clay were deposited in the river valley. The Evergreen Manor site is

located in the preglacial Rock River buried valley.

Overburden Geology

The surficial geology in the vicinity of the Evergreen Manor site consists of windblown sand and

silt, lacustrine sand, silts and clays, and outwash sand and gravel deposited within the preglacial

Rock River valley. Till deposits are found primarily along the valley margins. The valley was

CH01\PUBLIC\WO\RAC\036\29672S-3.WPD

RFW036-2A-AHVH

Section: 3 Revision: 0

Date: 31 October 2000

Page: 4 of 7

primarily filled with deposits from the Quaternary Period, during the Illinoian and Wisconsinan

glacial events. The sand and gravel deposits are the most abundant and most extensive deposits in

the buried valley, and can reach a thickness of up to 300 ft. in the vicinity of the Evergreen Manor

site (IDENR, 1960).

Bedrock Geology

The bedrock geology in the vicinity of the Evergreen Manor site is characterized by the Ordovician

and Cambrian clastic and carbonate rocks. The Galena and Platteville dolomite, and the Ancell

Groups represent the Ordovician Period in this area. The Ancell Group consists of two formations:

the Glenwood, which is a sandy shale, and the St. Peter, which is predominantly a well-sorted

sandstone up to 400 ft thick. The ancient Rock River eroded the Galena and Platteville dolomite,

and the Glenwood, and carved its valley into the St. Peter sandstone (Colten, 1986).

The Cambrian rocks are dominated by sandstones with lesser thicknesses of shale and dolomite. The

Potosi (dolomite) and Franconia (sandy shale) Formations separate the Ironton-Galesville sandstone

from the Ordovician rocks. The Ironton-Galesville sandstone has a thickness of up to 170 ft. The

Eau Claire Formation is up to 450 ft thick, and the Mt. Simon sandstone can be up to 1600 ft thick.

The sedimentary bedrock units in the vicinity of the Evergreen Manor site were deposited on an

irregular surface of metamorphic and igneous Precambrian rocks. Beneath the site, the Precambrian

consists of a granite (Colten, 1986).

3.6 SOILS

The predominant surficial soil type mapped for the site and surrounding area is the Warsaw loam

(USDA-NRCS, 1980).

CH01\PUBLIC\WO\RAC\036\29672S-3.WPD

RFW036-2A-AHVH

Section: 3 Revision: 0

Date: 31 October 2000

Page: 5 of 7

The Warsaw loam is a nearly level to gently sloping soil found on terraces, convex ridges, outwash

plains, gravelly kames and stream terraces. Depending on the slope, the surface layer is about 10

to 12 inches thick and consists of a very dark gray to a very dark brown loam. The subsoil is from

about 24 to 41 inches thick and consists of loam to gravelly loam and varies in color from dark

grayish brown to brown to dark reddish brown. The substratum, to a depth of about 60 inches,

consists of yellowish brown, calcareous sand and gravel. The permeability of the Warsaw loam is

moderate to rapid, with moderate water capacity, and moderate organic matter content (USDA-

NRCS, 1980).

Other soil types exist within the site area. Soil types located near the Rock River and Dry Creek are

characterized by higher clay contents and moderate permeabilities. Other soil types, further from

the water ways, are characterized by higher sand or sand and gravel contents and rapid permeability

(USDA-NRCS, 1980).

3.7 REGIONAL HYDROGEOLOGY

The sand and gravel deposits have significant permeability and transmissivity and are the

predominant local water source in the vicinity of the preglacial valley.

Hydrogeology

The unconsolidated outwash sand and gravel, the St. Peter, Ironton-Galesville, and Mt. Simon

Formation sandstones are the aquifers underlying the site. The outwash sands and gravels have

significant permeability and transmissivity and are the predominant local water source for private

residences in the vicinity of the preglacial Rock River Valley. Larger wells owned or used by

municipalities or developments draw groundwater from the bedrock aquifers.

CH01\PUBLIC\WO\RAC\036\29672S-3.WPD

RFW036-2A-AHVH

Section: 3 Revision: 0

Date: 31 October 2000

Page: 6 of 7

The outwash sand and gravel is an unconfined aquifer with more uniform (i.e. better sorted) deposits

at depth. Hydraulic conductivity in the outwash deposits has been estimated at 1x10<sup>-3</sup> cm/sec (IEPA,

1997). According to well logs, most of the private residential water wells are finished in the outwash

sand and gravel deposits at depths of 50 to 80 ft below ground surface (bgs). The water table is

approximately 35 ft bgs (U.S. EPA, 1999a).

The Galena and Platteville dolomite is an aquitard with a hydraulic conductivity estimated to range

from 1x10<sup>-8</sup> to 1x10<sup>-11</sup> cm/sec (IEPA, 1997). Flow in the dolomite is through vertically oriented

fracture and joint systems. Although not used for high yield production wells the Galena and

Platteville dolomite is used for small-demand, rural domestic and livestock water supply (Colten,

1986).

The Glenwood Formation, consisting of shale overlying a poorly sorted sandstone, has an estimated

hydraulic conductivity of 1x10<sup>-4</sup> to 1x10<sup>-7</sup> cm/sec (IEPA, 1997).

The St. Peter sandstone underlies the Galena and Platteville, and Glenwood Formations, except

along the axis of the Rock River buried valley, where the overlying bedrock has been removed by

erosion. The St. Peter sandstone has an estimated hydraulic conductivity of 1x10<sup>-4</sup> cm/sec and is

widely used as a water source in Winnebago County (IEPA, 1997).

The Ironton-Galesville Formation sandstone is a confined unit and a very productive aquifer.

However, due to its depth, few wells are finished in the Ironton-Galesville Formation (Colten, 1986).

The same holds true for the Mt. Simon sandstone, which can reach a thickness of 1600 ft, and

overlies the Precambrian granite.

CH01\PUBLIC\WO\RAC\036\29672S-3.WPD

RFW036-2A-AHVH

Section: 3 Revision: 0

Date: 31 October 2000

Page: 7 of 7

3.8 REGIONAL SURFACE WATER HYDROLOGY

The Evergreen Manor site lies in the Lower Rock River drainage basin. The Rock River orginates

in Wisconsin and enters Illinois south of Beloit. In Illinois, the Rock flows in a southwesterly

direction to its confluence with the Mississippi at Rock Island. In the vicinity of the site, the Rock

River flows generally in a north to south direction. The only tributary to the Rock on the site is Dry

Creek, which flows from the northeast to the southwest and discharges into the Rock in the

Hononegah Forest Preserve. Two lakes, Pearl Lake and Victoria Lake, are located north of the site,

west of Illinois Route 251. These lakes are former sand and gravel quarries.

3.9 GROUNDWATER USE IN THE AREA

The City of Roscoe is part of the North Park Public Water District (NPPWD), however, not all

residences receive their water from this source. At the time of this writing, most of the residential

wells within the Evergreen Manor site plume have been abandoned and these residences have been

connected to the public water supply. However, private wells are still in use on either side of the

plume and draw groundwater from the shallow sand and gravel aquifer. Two municipal wells

providing a portion of the water to the NPPWD are located at the corner of Hononegah Road and

Cedar Brook Road. These wells draw water from a depth of 750 feet below grade from the St. Peter

Sandstone (IEPA, 1997).

CH01\PUBLIC\WO\RAC\036\29672S-3.WPD

RFW036-2A-AHVH

Section: 4 Revision: 1 Date: 28 March 2001 Page: 1 of 11

#### **SECTION 4**

#### **ENVIRONMENTAL INVESTIGATION PROCEDURES**

This section describes the procedures used during the field investigation. The field investigation was conducted between 10 May 2000 and 7 June 2000 and consisted of the following activities:

- Fracture Trace Analysis
- CPT Groundwater Sampling
- Monitoring Well Sampling
- Residential Well Sampling
- Sediment Sampling
- Surface Water Sampling
- Groundwater Elevation Measurements
- Ecological Investigation

Prior to starting field activities, a site-specific Health and Safety Plan (HASP) (WESTON, 2000C), and a QAPP/FSP (WESTON, 2000a) were prepared. The HASP describes the safety protocols for field activities. The HASP was prepared in accordance with the Occupational Safety and Health Administration (OSHA) requirements as outlined in 29 CFR 1910 and other applicable requirements. The QAPP/FSP presents the organization, objectives, functional activities, and specific Quality Assurance and Quality Control (QA/QC) activities associated with the field activities. The QAPP/FSP also describes the specific protocols for sampling, sample handling and storage, chain of custody, and laboratory and field analysis. The QAPP/FSP was prepared in accordance with U.S. EPA QAPP guidance documents; in particular, the *Interim Guidelines and Specifications for Preparing Quality Assurance Project Plans* (QAMS-005/80) (U.S. EPA, 1980), *Region V Content Requirements for QAPPs* (U.S. EPA, 1989b), and the *Region V Model QAPP* (U.S. EPA, 1991d).

The laboratory analysis was performed by four different laboratories. On-site VOC analysis of soil and water samples was performed by Lockheed Martin Services Group, Environmental Services & Technologies Region 5 (ESAT), of Chicago, Illinois. Off-site VOC analysis of water samples was

Section: 4 Revision: 1

Date: 28 March 2001 Page: 2 of 11

performed by Mitkem Corporation, of Warwick, Rhode Island (MITKEM). Off-site VOC analysis

of soil samples was performed by Datachem Laboratories of Salt Lake City, Utah (DATACHEM).

Off-site water quality parameters analyses were performed by Chemtech, of Edison, New Jersey

(CHEMTECH).

4.1 FRACTURE TRACE ANALYSIS

A fracture trace analysis was performed in order to better identify potential source areas, potential

locations for CPT work, and to obtain a better understanding of the flow system. Resolution

Resources, Inc. (RRI), of Minneapolis, Minnesota and Warrenton, Virginia were contracted to

perform this study. RRI used aerial photographs dating back to 1939 in order to identify linear

features in the overburden (lineaments), that may have been caused by fractures in the underlying

bedrock. This information was used to target potential preferential flow pathways for subsequent

groundwater profile sampling and analysis. A complete discussion of this approach is presented in

RRI's report, which is included in Appendix B.

4.2 CONE PENETROMETER TESTING (CPT)

A 20 ton, truck-mounted cone penetrometer rig was used to perform cone penetrometer testing

(CPT), and groundwater sampling between 25 May and 6 June 2000. The purpose of CPT was to

identify the stratigraphy at various locations across the site, and then use that information to choose

groundwater sampling depth intervals at each location. Of the 13 locations identified with "CPT",

on Figure 4-1, stratigraphy testing was performed at 10 of them.

The stratigraphy at each CPT location was analyzed by hydraulically pushing a 2-inch diameter,

instrumented probe into the subsurface and recording geotechnical data. This included end-bearing

resistance, friction along the sides of the probe, and electrical conductivity. End bearing resistance,

measured in tons per square foot (tsf), helps differentiate between different geologic materials. For

CHLAN01\WP\WO\RAC\036\29672S-4.WPD

RFW036-2A-AHVH

Section: 4 Revision: 1

Date: 28 March 2001

Page: 3 of 11

instance, sand has a greater end bearing resistance than a clay. Similarly, the amount of friction

along the sides of the probe helps identify the type of geologic material. Lower friction is incurred

when pushing the probe through a sand than when pushing the probe through a clay, silty clay, or

silt. The ratio of friction to end bearing resistance (friction ratio) is a measure that helps identify the

amount of fine grained material (clay and silt) present. As an example, when pushing through a

predominantly silty sand, both the end bearing resistance and the friction ratio will be high.

The electrical conductivity measure was used because of its ability to identify features such as the

water table, groundwater plumes, and clay lenses. Zones that are unsaturated, or saturated with non-

conductive liquids, will give low electrical conductivity readings; saturated zones will give

significantly higher readings. This measure is also useful in identifying dissolved constituents in

groundwater plumes, however, this was not an expectation at the Evergreen Manor site.

CPT was conducted at locations CPT01 through CPT08, CPT11, and CPT13. As described in the

Field Sampling Plan, because of the expected similarity in stratigraphy across the site, CPT was not

intended to be conducted at each location. Therefore, CPT was not conducted at CPT09, CPT10,

and CPT12.

The expectation was to advance CPT holes to a depth of about 120 feet below grade, however, the

abundance of gravels and cobbles limited the depth of penetration. In most instances, the CPT holes

were advanced to at least 90 feet below grade. However, at CPT07, CPT08, and CPT13, refusal was

encountered at depths less than 10 to 15 feet below grade. It was noted that a concrete-containing

fill may be present in this area, which could cause refusal of the CPT probe. It is also possible that

natural features (cobbles and/or boulders) were responsible for the refusals. Several attempts were

made to penetrate deeper at each of these locations without success.

CHLAN01\WP\WO\RAC\036\29672S-4.WPD

RFW036-2A-AHVH

Section: 4 Revision: 1

Date: 28 March 2001 Page: 4 of 11

4.3 CPT GROUNDWATER SAMPLING

Groundwater samples were collected between 25 May and 6 June 2000. In general, groundwater

samples were collected directly after completion of the CPT hole (stratigraphy hole). Based on the

stratigraphy, sampling intervals were chosen. The groundwater sampling locations were typically

within several feet of the stratigraphy hole locations. In the same manner that the CPT rods were

pushed, the groundwater sampler was advanced by hydraulically pushing it to the pre-determined

depth. The groundwater sampler consisted of a screen with a retractable outer casing. The screen

was opened by pulling back on the rods, exposing the screen to the native soil and groundwater.

At CPT-01, the groundwater sampler was first pushed to the deepest location to collect a

groundwater sample. The sampler was then pulled back to the next shallower depth to collect

another sample. This was repeated until the shallowest sample was collected. Difficulties were

encountered while attempting to purge the rod string of water. Because of the inefficiency of

purging the rod string, this approach was changed for all subsequent CPT groundwater samples;

instead of this bottom-up approach, it was changed to a top-down approach.

For locations CPT-02 through CPT-13, the groundwater sampler was first pushed to the shallowest

depth interval to collect a sample. The entire rod string was then pulled out, decontaminated, and

re-deployed into the same hole to the next sampling interval. In some instances, only a stratigraphy

hole and one other hole were pushed. However, at CPT-03 the sampling hole was reamed out in the

vadose zone, from multiple deployments through the same hole, to the point where the hole did not

provide enough lateral support to the rod string. The result was that the rod string broke just below

the water table in CPT-03. Subsequently, several sampling holes were typically pushed at each

location.

At each sampling depth the groundwater sampler was opened, and small diameter tubing, with a

check valve at the bottom was lowered through the CPT rods. Groundwater was pumped from the

CHLAN01\WP\WO\RAC\036\29672S-4.WPD

RFW036-2A-AHVH

Section: 4
Revision: 1

Date: 28 March 2001 Page: 5 of 11

tubing in conjunction with a peristaltic pump, however, groundwater could not be pumped to the

surface due to the depth of the water table. To collect the groundwater samples, the tubing was

pulled from the rod string and evacuated using the peristaltic pump. At the time of sampling,

measurements of pH, specific conductance, temperature, turbidity, DO, ORP, and Fe<sup>+2</sup> were recorded

on Water Sample Collection Forms, which are presented in Appendix C. Care was taken to fill the

40 mL VOA vials at an angle to minimize splashing and bubbling, and to ensure that they were

closed with no headspace. All of the samples collected were analyzed for VOCs.

4.4 MONITORING WELL SAMPLING

A total of 15 existing monitoring wells, shown on Figure 4-1, were sampled from 30 May through

2 June 2000. Note: The monitoring wells are identified with a prefix of 'G' (e.g. G103D) on the

figures, which was the designation given by the IEPA at installation time. However, they are

referred to throughout the text with the prefix 'MW' (e.g. MW103D), which corresponds to the

designated sample numbering system in the approved QAPP.

Each sample collected from the monitoring wells were analyzed for VOCs and Fe<sup>+2</sup> (in the field).

Additional volume was collected from the deep well at each cluster and from non-clustered wells

for water quality parameter analysis. Water quality parameters were analyzed to aid in evaluating

the potential for biodegradation of the chlorinated hydrocarbon plume. Water quality parameters

include: ammonia, chemical oxygen demand (COD), nitrate, nitrite, orthophosphate, sulfate, sulfide,

and ferrous iron (Fe<sup>2+</sup>; measured in the field).

Each monitoring well, with the exception of MW-112, was purged and sampled using a

decontaminated Grundfos<sup>TM</sup> pump and tubing. Monitoring well MW-112 was found to be bent and

the pump could not fit down the stainless steel riser. A disposable bailer was cut to a length of 7

inches in order to fit past the bent riser, and was used to purge and sample MW-112. The depth to

water in the well and the total depth of the well were measured with an electrical sounding device.

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RFW036-2A-AHVH

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Section: 4 Revision: 1

Date: 28 March 2001

Page: 6 of 11

The top of the inner well casing was used as the reference point for these measurements. These

measurements were used to calculate well volume and were recorded, along with the time, on Water

Sample Collection Forms, which are presented in Appendix C. A minimum of three purge volumes

was removed from the wells. After removing the third well volume, field measurements of pH,

specific conductance, temperature, and turbidity were recorded.

Purging continued until the measurements for all parameters had stabilized ( $\pm 0.25$  units for pH,  $\pm 10$ 

percent for specific conductance,  $\pm 1.0$  C, and  $\pm 10$  percent for turbidity) for two consecutive rounds

of readings or until five well volumes had been purged. Measurements of dissolved oxygen (DO),

and oxygen reduction potential (ORP) were recorded, but were not used to determine stabilization.

A Hach Test Kit (Model IR-18C) was used to measure Fe<sup>+2</sup> and samples were collected once the well

had stabilized. Sample containers (40 mL VOA vials and three 1 L plastic bottles) were filled

directly from the pump tubing. Care was taken to fill the VOA vials at an angle to minimize

splashing and bubbling, and to ensure that they were closed with no headspace.

4.5 RESIDENTIAL WELL SAMPLING

A total of 22 residential wells, shown on Figure 4-2, were sampled between 31 May and 6 June

2000. All of the samples collected were analyzed for VOCs and one water quality parameter (Fe<sup>+2</sup>).

Six of the 22 samples were additionally analyzed for the following water quality parameters:

ammonia, COD, nitrate, nitrite, orthophosphate, sulfate, and sulfide. Water quality parameters were

analyzed to aid in evaluating the potential for biodegradation of the chlorinated hydrocarbon plume.

Each residential well was sampled at an outside cold-water spigot. To purge each residential well.

water was allowed to flow for at least 20 minutes. After the 20 minutes, field measurements of pH.

specific conductance, temperature, and turbidity were recorded. Purging continued until two

consecutive rounds of parameter measurements had stabilized (±0.25 units for pH, ±10 percent for

specific conductance, ±1.0 C, and ±10 percent for turbidity). DO and ORP measurements were

CHLAN01\WP\WO\RAC\036\29672S-4.WPD

RFW036-2A-AHVH

Section: 4 Revision: 1

Date: 28 March 2001

Page: 7 of 11

taken, however, they were not used to determine stabilization. All measurements were recorded on

Water Sample Collection Forms, which are presented in Appendix C.

A Hach Test Kit (Model IR-18C) was used to measure Fe<sup>+2</sup> and samples were collected once the well

had stabilized. Sample containers, 40 mL VOA vials and three 1 L plastic bottles glass (as

applicable for water quality parameters), were filled directly from the spigot, by first filling the VOA

vials, and then the plastic bottles. Care was taken to fill the VOA vials at an angle to minimize

splashing and bubbling, and to ensure that no headspace remained once a vial was capped.

4.6 <u>SEDIMENT SAMPLING</u>

A total of 6 sediment samples and one QC field duplicate were collected from Dry Creek and the

Rock River on 23 and 24 May 2000, at locations shown on Figure 4-3. The sediment samples were

collected to determine if site contaminants were present in river sediments. Three sediment samples

(SD-01 to SD-03) were collected at regular intervals along the eastern bank of the Rock River. SD-

01 was an upriver location. SD-02 was upriver of where the plume was believed to discharge to the

Rock River, but downgradient of the confluence with Dry Creek. SD-03 was adjacent to the site.

The other three sediment samples (SD-04 to SD-06) were collected from Dry Creek. SD-04 was an

upgradient location. SD-05 and SD-06 were collected from locations within and downstream of the

site.

The sediment samples were collected by shoveling sediment from the river or creek and then placing

the sediment in a 16-ounce glass jar using a sterilized plastic scoop. Sediment sampling was

conducted from downriver to upriver locations, to minimize the impact of sediment disturbance

and/or cross contamination of samples. All reusable field sampling equipment used for sediment

sampling was decontaminated between sample locations. The six investigative sediment samples

and a QC duplicate were submitted to the mobile ESAT laboratory for VOC analysis and to

CHEMTEC for Total Organic Carbon (TOC) analysis. One sample (SD04-01) and its QC duplicate

CHLAN01\WP\WO\RAC\036\29672S-4.WPD

RFW036-2A-AHVH

Section: 4

Revision: 1 Date: 28 March 2001 Page: 8 of 11

(SD04-01) were shipped to DATACHEM for confirmation VOC analysis. The analytical results of

the sediment sampling are presented in Subsection 5.5.

4.7 SURFACE WATER SAMPLING

A total of six surface water samples (SW01-SW06) were collected from Dry Creek and the Rock

River on 23 and 24 May 2000, as shown on Figure 4-3. The surface water samples were collected

to determine if site contaminants were migrating into Dry Creek and the Rock River. Surface water

samples were collected at approximately the same locations as sediment samples. At all six

locations, surface water samples were collected prior to collection of sediment samples. Surface

water samples were collected from the bank directly into pre-preserved 40 mL VOA vials. The

sampling progressed from downstream to upstream locations to minimize the impact of sediment

disturbance and/or cross contamination of samples.

The six samples and one QC duplicate sample were submitted to the mobile ESAT laboratory for

VOC analysis. One sample (SW04-01) and its QC duplicate (SW04-01DP) were submitted to

MITKEM for confirmation VOC analysis. The analytical results of the surface water sampling are

presented in Subsection 5.6.

4.8 GROUNDWATER ELEVATION MEASUREMENTS

Depth to groundwater and depth to bottom were measured at 16 monitoring wells, on 6 June 2000,

over a period of about 1.5 hours. An electrical sounding device (Solinst water level indicator) was

used to measure the depth to water and the total depth of each well. The top of the inner well casing

was used as the reference point for these measurements. Although bent, the top of the well casing

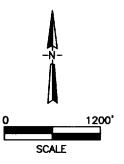
at MW-112 was still used as a reference point. This should impart an error of only several tenths

of a foot and since it is the furthest upgradient well location, the data is deemed usable.

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RFW036-2A-AHVH

ELEVATOR RD.



NOTE:
BASEMAP ADAPTED FROM FIGURE 1 OF THE CONTAMINANT SOURCE
EVALUATION — EVERGREEN MANOR SITE, DATED JANUARY 1997, BY
CONESTOGA—ROVERS & ASSOCIATES.

FIGURE 4-1

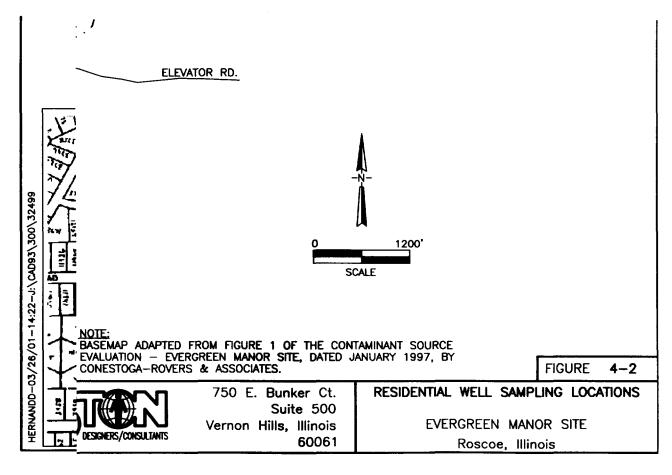


HERNANDD-03/26/01-14:21-J:\CAD93\200\23300

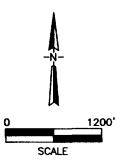
750 E. Bunker Ct. Suite 500 Vernon Hills, Illinois 60061 GROUNDWATER SAMPLING LOCATIONS
(CPT AND MONITORING WELLS)

EVERGREEN MANOR SITE

Roscoe, Illinois



ELEVATOR RD.



NOTE:
BASEMAP ADAPTED FROM FIGURE 1 OF THE CONTAMINANT SOURCE
EVALUATION — EVERGREEN MANOR SITE, DATED JANUARY 1997, BY
CONESTOGA—ROVERS & ASSOCIATES.

FIGURE 4-3



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750 E. Bunker Ct. Suite 500 Vernon Hills, Illinois 60061 SEDIMENT AND SURFACE WATER
SAMPLING LOCATIONS
EVERGREEN MANOR SITE
Roscoe, Illinois

Section: 5 Revision: 1

Date: 28 March 2001

Page: 1 of 35

**SECTION 5** 

**ENVIRONMENTAL INVESTIGATION RESULTS** 

This section presents results of data collected during the various environmental investigations. The

investigations included a fracture trace analysis, CPT groundwater sampling, monitoring well

sampling, residential well sampling, sediment sampling, surface water sampling, groundwater

elevation measurements, and ecological investigation.

5.1 FRACTURE TRACE ANALYSIS

The fracture trace analysis was completed by RRI according to the procedures outlined in Subsection

4.1. RRI analyzed a number of aerial photographs and photo pairs in order to identify lineaments

in the overburden, which could indicate the presence of underlying faults. The report by RRI is

included in Appendix B. The following summarizes the findings of the fracture trace analysis.

The fracture trace analysis identified two main sets of fracture trends: a north-south/east-west set,

and a northeast-northwest conjugate set. To the north of Hononegah Road, the predominant set of

fractures is the northeast-northwest set, and to the south of Hononegah Road, the predominant set

of fractures is the north-south/east-west set. This information supports the pre-RI plume location

(shown on Figure 4-1), derived from previous investigations, and also indicates that a predominant

flow path (or direction) may exist. Thus, the flow may be influenced by fractures propagated into

the overburden from the underlying bedrock.

RRI also indicated, that based on the predominant fracture set directions, a possible source area

could be identified as the industrial park near the intersection of Route 251 and Rockton Road. They

also identified another potential source, which is a former farm located near Hononegah Road and

upgradient of the residential area of the site.

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Section: 5

Revision: 1 Date: 28 March 2001

Page: 2 of 35

Based on the fracture trace analysis, RRI suggested over 40 points located along fractures, or at

fracture junctions, where CPT groundwater sampling could be done in order to better delineate the

plume, or identify a source area. As a result of this recommendation, all the CPT locations along

McCurry Road were adjusted to coincide with identified fracture locations, and two additional

locations (CPT 1 and CPT 2) were completed along Route 251 near the potential source area.

5.2 CPT GROUNDWATER SAMPLING

CPT groundwater samples were collected at 10 locations. At each of these locations a sample was

collected from a minimum seven different sampling depths. At two of the locations, samples were

collected from eight depths. Each sample was analyzed for VOCs by the on-site laboratory (ESAT)

as well as for Fe<sup>+2</sup>, which was measured in the field at the time the sample was collected. Other field

parameters included temperature, pH, specific conductance, turbidity, dissolved oxygen, and

oxidation-reduction potential.

A total of 72 CPT groundwater samples were collected. Additionally, six OC duplicate samples

were analyzed on site by ESAT, 10 confirmation samples were analyzed off site by MITKEM, and

two QC duplicate samples of the 10 off site analysis samples were analyzed by MITKEM. All of

these samples were analyzed for VOCs. Groundwater samples were not collected at locations

CPT07, CPT08, and CPT13 due to shallow refusal of the CPT sampling equipment. The procedures

used for the CPT groundwater sampling are described in Subsection 4.3, and the CPT groundwater

results for VOCs are presented in Table 5-1. The field parameter results are contained on the Water

Sample Collection Forms, attached in Appendix C. The analytical data are attached in Appendix D.

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Section: 5 Revision: 1

Date: 28 March 2001 Page: 3 of 35

#### **VOC Results**

Thirteen VOCs were detected above method detection limits in the groundwater samples. These detections are summarized below:

- Acetone was detected in 18 samples at concentrations from 2 μg/L in CPT12-05, CPT12-06, and CPT12-07, to 470 μg/L in CPT02-03.
- Methylene chloride was only detected in sample CPT03-05 at a concentration of 0.5 μg/L.
- 1,1-Dichloroethane was only detected in sample CPT11-05 and its duplicate at a concentration of 2 μg/L.
- 2-Butanone was only detected in sample CPT05-06 at a concentration of  $16 \mu g/L$ .
- Cis-1,2-dichloroethene was detected in five samples at a concentration of 1 μg/L in samples CPT01-03 through CPT01-06, and at a concentration of 2 μg/L in sample CPT01-02.
- 1,1,1-Trichloroethane was detected in 21 samples from locations CPT01, CPT03, CPT10 and CPT11. Detected concentrations ranged from 0.7 μg/L in sample CPT11-08 to 3 μg/L in samples CPT11-05, and CPT11-06.
- Benzene was only detected in sample CPT09-01 and duplicate sample CPT12-04DUP at a concentration of 0.5 μg/L, and in sample CPT09-07 at concentration of 0.6 μg/L.
- Trichloroethene was detected in all of the 8 samples from location CPT01 at concentrations from 2 to 4 μg/L.
- Toluene was detected in 75 samples, and at least once at each CPT sampling location.

  Detected concentrations ranged from 0.5 μg/L to 3 μg/L.

Section: 5 Revision: 1

Date: 28 March 2001

Page: 4 of 35

Tetrachloroethene was only detected in sample CPT10-04 and its duplicate at a

concentration of 0.6  $\mu$ g/L, and in sample CPT10-02 at a concentration of 0.9  $\mu$ g/L.

Ethylbenzene was only detected in sample CPT09-07 at a concentration of  $0.6 \mu g/L$ .

• Xylenes were detected at concentrations of 0.5 μg/L, in sample CPT02-02, 0.6 μg/L

in samples CPT02-01, CPT02-07, and CPT11-01, and 0.7 μg/L in samples CPT06-01

and CPT09-07.

Field Parameter Results

• DO was detected at concentrations from 5.2 to 12.0 mg/L. Most of the results were

found to be between 8 and 10 mg/L.

• ORP was measured at concentrations from -199 to 155 mV. These values were fairly

evenly distributed between the maximum and minimum and were not found to be

dependent on depth.

Fe<sup>+2</sup> was detected in 17 samples up to a concentration of 1.1 mg/L.

5.3 MONITORING WELL SAMPLING

Groundwater samples were collected from 15 monitoring wells and analyzed for VOCs. Eight of

the samples, primarily from the deep wells, were also analyzed for water quality parameters. The

water quality parameters included ammonia, COD, nitrate, nitrite, orthophosphate, sulfate, sulfide,

and Fe<sup>+2</sup>. Each sample was analyzed in the field for dissolved oxygen, oxidation-reduction potential,

and Fe<sup>+2</sup>. The procedures used for monitoring well sampling are described in Subsection 4.4. The

VOC and water quality parameter results are presented in Tables 5-2 and 5-3. Field parameters are

presented on Table 5-3 for samples which were also analyzed for water quality parameters. Field

parameters for the remaining samples are found on the Water Sample Collection Forms attached in

Appendix C. Analytical data are attached in Appendix D.

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Section: 5 Revision: 1

Date: 28 March 2001 Page: 5 of 35

**VOC Results** 

Six VOCs were detected above method detection limits in the monitoring well samples and are listed

below:

• 1,1,2-Trichloro-1,2,2-trifluoromethane was detected in sample MW103S at a

concentration of 2 µg/L, and in sample MW103D and its QC duplicate at

concentrations of 300 and 180 µg/L, respectively.

• Cis-1,2-Dichloroethene was detected in sample MW105S at a concentration of 1

 $\mu$ g/L, and in sample MW105D and its QC duplicate at concentrations of 1  $\mu$ g/L and

2 μg/L, respectively.

• 1,1,1-Trichloroethane was detected in eight samples at concentrations from 1 to 3

μg/L from MW103D, MW104D, MW104S, MW105D, MW105S, and MW112.

• Trichloroethene was detected in four samples at concentration from 2 to 3  $\mu$ g/L, from

MW101D, MW105D and MW105S.

Tetrachloroethene was detected in eight samples at concentrations from 0.5 to 9

μg/L, from MW103D, MW103S, MW105D, MW105S and MW108D.

Water Quality Parameters

The results of the water quality analyses for monitoring well samples are presented in Table 5-3, and

are summarized below:

Ammonia and sulfide concentrations in all wells were below their minimum

detectable concentrations.

• COD was detected in well MW112 at a concentration of 10 mg/L.

• Nitrite was detected in MW101D at a concentration of 7.8 mg/L.

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Section: 5 Revision: 1

Date: 28 March 2001 Page: 6 of 35

• Nitrate concentrations range from 3.3 mg/L in well MW112 to 8.5 mg/L in MW108D.

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Orthophosphate concentrations range from non-detectable in MW102D, MW105D

and MW108D to 0.69 mg/L in MW103D.

• Sulfate concentrations range from 15 mg/L in MW102D to 27 mg/L in MW101D.

Field Parameter Results

• DO was detected at concentrations from 2.8 to 7.7 mg/L. Most of the readings were

between 5.5 and 8 mg/L, however, readings from MW104S, MW105S, and

MW105D were 3.2, 4.8, and 2.8 mg/L, respectively.

• ORP was measured from 4 to 213 mV. Most of these measurements were found to

be greater than 100 mV.

• Fe<sup>+2</sup> was not detected in any of the monitoring well samples.

5.4 RESIDENTIAL WELL SAMPLING

Groundwater samples were collected from 22 residential wells. Each sample was analyzed for

VOCs by the on-site laboratory (ESAT) as well as for Fe<sup>+2</sup>, which was measured in the field at the

time the sample was collected. Of the 22 total samples collected, two were also sent to an off-site

laboratory for confirmation VOC analysis. At locations RW01 through RW05 and RW07,

additional volume was collected for water quality parameter analysis, which included ammonia.

COD, nitrate, nitrite, orthophosphate, sulfate, and sulfide. The procedures for residential well

sampling are described in Subsection 4.5. The VOC and water quality parameter results are

presented in Tables 5-4 and 5-5. Field parameters are presented on Table 5-5 for samples also

analyzed for water quality parameters. Field parameters for the remaining samples are found on the

sample collection forms attached in Appendix C. Analytical data are attached in Appendix D.

CHLAN01\WP\WO\RAC\036\29672S-5.WPD

RFW036-2A-AHVH

Section: 5 Revision: 1 Date: 28 March 2001

Page: 7 of 35

#### **VOC Results**

Six VOCs were detected above method detection limits in the residential well samples and are listed below:

- Acetone was detected in sample RW03 and duplicate sample RW05 at concentrations of 0.8 μg/L and 0.6 μg/L, respectively.
- Chloroform was detected only in sample RW08 at a concentration of 0.9 μg/L.
- Cis-1,2-Dichloroethene was detected in samples from RW04 at a concentrations of 1 and 2 μg/L.
- 1,1,1-Trichloroethane was detected in 11 samples from RW03, RW04, RW07, RW08, RW11 and RW19 at concentrations from 1 to 5 μg/L.
- Trichloroethene was detected in five samples from RW04, RW07 and RW19 at concentrations from 0.7 to 6 µg/L.
- Tetrachloroethene was detected at a concentration of 2 μg/L in samples from RW04, and at a concentration of 0.9 μg/L in sample RW19 and its duplicate.

#### **Water Quality Parameters**

The results of the water quality analyses for residential well samples are presented in Table 5-5.

- COD, ammonia, and nitrite were not detected in any of the residential well samples above their respective method detection limits.
- Nitrate was detected in the samples from RW02 and RW04 at concentrations of 3.9 and 6.3 mg/L, respectively.
- Orthophosphate was detected in samples from RW01, RW03, and RW07.
   Concentrations detected ranged from 0.029 mg/L in the sample from RW07 to 0.051 mg/L in the samples from RW01 and RW03.

Revision: 1

Date: 28 March 2001

Page: 8 of 35

• Sulfate was detected in samples from RW04 and RW07 at concentrations of 19 and

28 mg/L, respectively.

Sulfide was detected in samples from RW03 and RW04 at concentrations of 2.4 and

1.6 mg/L, respectively.

Field Parameter Results

• DO concentrations were found from 3.25 to 9.2 mg/L in the residential wells. Most

of the measurements were found to be below 6 mg/L.

• ORP was measured from 80 to 238 mV in the residential wells. Measurements were

fairly well distributed between these values.

• Fe<sup>+2</sup> was not detected in any of the residential well samples.

5.5 SEDIMENT SAMPLING

A total of 6 sediment samples (SD01 through SD06) were collected during the field investigation

from Dry Creek and the Rock River and analyzed for VOCs and TOC. One sample and a duplicate

from SD04 were sent to an off-site laboratory (DATACHEM) for VOC and TOC analysis. The

procedures used for the sediment sampling are described in Subsection 4.6. The sediment sample

results are presented in Tables 5-6 and 5-7 and the analytical data are attached in Appendix D.

• Benzene was detected only in the duplicate sample analyzed by DATACHEM, from

location SD04, at a concentration of 2 ug/kg.

• Chloroform was detected only in the sample from location SD01 at a concentration

of 8 ug/kg.

Methyl acetate (Methylene chloride) was detected in samples from SD01, SD02, and

SD05 at concentrations of 5, 9, and 9 ug/kg, respectively.

Toluene was detected only in the duplicate sample analyzed by DATACHEM, from

location SD04, at a concentration of 0.7 ug/kg.

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Section: 5 Revision: 1

Date: 28 March 2001

Page: 9 of 35

• TOC was detected in all of the samples at concentrations from 4,200 mg/kg in the

sample from SD04 to 30,000 mg/kg in the sample from SD03.

5.6 SURFACE WATER SAMPLING

Six surface water samples were collected from Dry Creek and the Rock River at approximately the

same locations as the sediment samples. Each sample was analyzed for VOCs. One sample and a

QC duplicate from SW04 were sent to an off-site laboratory (MITKEM) for confirmation VOC

analysis. The procedures used for surface water sampling are described in Subsection 4.7. The

surface water sample results are presented in Table 5-8 and the analytical data are attached in

Appendix D.

VOCs were not detected in any of the surface water samples.

5.7 GROUNDWATER ELEVATION MEASUREMENTS

The depth to groundwater was measured in 16 monitoring wells, as described in section 4.8. The

depth to groundwater measurement and the elevation of the reference point were used to calculate

the elevation of groundwater in each well. Groundwater elevations were from 722.58 to 735.22 feet

above MSL. In each of the well clusters, the groundwater elevation in the shallow and deep wells

was found to be nearly identical. The biggest difference in groundwater elevations at any well

cluster measured was 0.08 ft.

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Sample ID:	CPT-01-01	CPT-01-02	CPT-01-03	CPT-01-04	CPT-01-05	CPT-01-06	CPT-01-07	CPT-01-08
Sample Number:								
Sample Date:	5/25/00	5/25/00	5/25/00	5/25/00	5/25/00	5/25/00	5/25/00	5/25/00
Laboratory:	ESAT							
Sample Screen Depth (ft below ground):	99	89	79	69	59	49	39	29
Units:	ug/L							
Parameter:								
Dichlorodifluoromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 υ	1 U
chloromethane	1 0	1 U	1 U	1 0	1 U	1 U	1 0	1 U
Vinyl Chloride	1 0	1 0	1 0	1 0	1 U	1 0	1 0	1 U
Bromomethane	1 0	1 U	1 U	1 0	1 U	1 U	1 U	1 U
Chioroethane	1 0	1 0	1 U	1 0	1 U	1 U	1 0	1 U
Trichloroffuoromethane	1 0	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1,2-Trichioro-1,2,2-trifluoromethane	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
1.1-Dichloroethene	1 U	1 U	1 U	1 Ü	1 U	1 U	1 U	1 U
Acetone	2 UJ	15 B	2 UJ					
Carbon Disulfide	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Methylene Chloride	1 0	1 U	1 U	1 0	1 U	1 U	1 U	1 U
Methyl Acetate	1 0	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trans-1,2-Dichloroethene	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
1,1-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Butanone	11 U							
Cis-1,2-Dichloroethene	4 U	2 J	1 J	1 J	1 J	1 J	4 U	4 U
Methyl tert-Butyl Ether	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1.2-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroform	3 U	3 U	3 U	3 U	3 U	3 U	3 U	3 U
Carbon Tetrachloride	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
1,1,1-Trichloroethane	2 U	1 J	1 J	1 J	1 J	2 U	1 J	2 U
Benzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	2 J	3 J	4 J	4 J	3 J	4 J	3 J	2 J
1,2-Dichioropropane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromodichloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Cis-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trans-1,3-Dichioropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1,2-Trichioroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Dibromochioromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	1 Ü	1 U	1 U	1 U	1 U	1 U	1 U	1 U
4-Methyl-2-pentanone	5 U	5 U	5 U	5 U	5 U	5 U	5 Ü	5 U
Toluene	0.7 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Tetrachloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Hexanone	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U
1,2-Dibromoethane	1 U	1 U	1 U	1 U	1 U	1 U	1 0	1 U
Chlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Ethylbenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-&/or p-Xylene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
o-xylene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Xylenes (total)								
Styrene	3 U	3 U	3 U	3 U	3 U	3 U	3 U	3 U
1,1,2,2-Tetrachioroethane	1 UJ							
1,3-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U_	1 U
1,4-Dichlorobenzene	1 υ	1 U	1 U	1 U	1 U	1 υ	1 U	1 υ
1,2-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2,4-Trichlorobenzene								
1,2-Dibromo-3-chloropropane						***		
Bromochloromethane								

Sample ID:	CPT-02-01	CPT02-01	CPT-02-02	CPT-02-03	CPT02-03	CPT-02-04	CPT-02-05	CPT-02-06
Sample Number:		EABX3	_		EABX4			
Sample Date:	6/2/00	8/2/90	6/2/00	6/2/00	6/2/00	\$/2/00	6/2/00	6/2/00
Laboratory:	ESAT	MitKem	ESAT	ESAT	MitKem	ESAT	ESAT	ESAT
Sample Screen Depth (ft below ground):	28	28	34	42	42	51	68	78
Units:	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Parameter:						<u> </u>		<u> </u>
Dichlorodifluoromethane	1 U		1 U	1 U	_	1 U	1 U	1 U
chioromethane	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
Vinyl Chloride	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
Bromomethane	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
Chloroethane	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
Trichlorofluoromethane	1 U		1 U	1 U		1 U	1 U	1 U
1.1.2-Trichloro-1.2.2-trifluoromethane	2 UJ		2 UJ	2 UJ		2 UJ	2 UJ	2 UJ
1,1-Dichloroethene	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
Acetone	2 UJ	5 U	2 UJ	55 J	470	2 UJ	2 UJ	2 UJ
Carbon Disulfide	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
Methylene Chloride	1 U	2 U	1 Ü	1 U	10 U	1 U	1 U	1 U
Methyl Acetate	1 UJ	***	1 UJ	1 UJ		1 UJ	1 UJ	1 UJ
Trans-1,2-Dichloroethene	2 U	1 U	2 U	2 U	5 U	2 U	2 U	2 U
1,1-Dichloroethane	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
2-Butanone	11 UJ	5 U	11 UJ	11 UJ	25 U	11 UJ	11 UJ	11 UJ
Cis-1.2-Dichloroethene	4 U	1 U	4 U	4 U	5 U	4 U	4 U	4 U
Methyl tert-Butyl Ether	1 UJ	_	1 UJ	1 UJ		1 UJ	1 UJ	1 UJ
1.2-Dichloroethane	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 υ
Chloroform	3 U	1 Ü	3 U	3 U	5 U	3 U	3 U	3 U
Carbon Tetrachloride	2 U	1 U	2 U	2 U	5 U	2 U	2 U	2 U
1,1,1-Trichloroethane	2 U	1 U	2 U	2 U	5 U	2 U	2 U	2 U
Benzene	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
Trichioroethene	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
1,2-Dichloropropane	1 U	1 U	1 Ü	1 U	5 Ü	1 U	1 U	1 U
Bromodichioromethane	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
Cis-1,3-Dichioropropene	1 U	1 Ü	1 U	1 U	5 U	1 U	1 U	1 U
Trans-1,3-Dichloropropene	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
1,1,2-Trichloroethane	1 U	1 0	1 U	1 U	5 U	1 U	1 U	1 U
Dibromochloromethane	1 U	1 1	1 U	1 U	5 U	1 U	1 U	1 U
Bromoform	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
4-Methyl-2-pentanone	5 U	5 U	5 U	5 U	25 U	5 U	5 U	5 U
Toluene	0.5 J	0.5 J	0.6 J	0.7 J	5 U	0.5 J	1 J	1 J
Tetrachloroethene	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
2-Hexanone	5 U	5 U	5 U	5 U	25 U	5 U	5 U	5 U
1,2-Dibromoethane	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
Chiorobenzene	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
Ethylbenzene	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
m-8/or p-Xylene	1 U		0.5 J	1 U		1 U	1 U	1 U
o-xylene	1 U		1 U	1 U		1 U	1 U	1 U
Xylenes (total)		0.6 J			5 U			
Styrene	3 UJ	1 U	3 00	3 UJ	5 U	3 UJ	3 UJ	3 UJ
1,1,2,2-Tetrachloroethane	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
1,3-Dichlorobenzene	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
1,4-Dichlorobenzene	1 U	1 0	1 U	1 U	5 U	1 U	1 Ü	1 U
1,2-Dichlorobenzene	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U
1,2,4-Trichlorobenzene	<u> </u>	1 U			5 U			
1,2-Dibromo-3-chloropropane		1 R			5 R			
Bromochloromethane		1 U		<u> </u>	5 U			

Sample ID:	CPT-02-07	CPT-03-01	CPT-03-02	CPT-03-02DP	CPT-03-03	CPT-03-04	CPT-03-05	CPT-03-06
Sample Number:								
Sample Date:	8/2/00	5/26/00	5/26/00	5/26/00	5/26/00	5/26/00	5/26/00	\$/2/00
Laboratory:	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT
Sample Screen Depth (ft below ground):	84	29	42	42	54		84	92
Units:	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Parameter:								
Dichlorodifluoromethane	1 0	1 U	1 U	1 U	1 U	1 U	1 U	1 U
chloromethane	1 0	1 U	1 U	1 U	1 U	1 U	1 U	1 0
Vinyl Chloride	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromomethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 0
Chloroethane	1 U	1 Ü	1 U	1 U	1 U	1 U	1 U	1 U
Trichlorofluoromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1,2-Trichloro-1,2,2-trifluoromethane	2 UJ	2 U	2 U	2 U	2 U	2 U	2 UJ	2 UJ
1,1-Dichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Acetone	2 UJ	20 B	32 B	19 B	20 B	25 B	9 J	2 UJ
Carbon Disulfide	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Methylene Chloride	1 υ	1 U	1 U	1 U	1 U	1 U	0.5 J	1 U
Methyl Acetate	ט 1 ט	1 U	1 U	1 U	1 Ü	1 U	1 UJ	1 ບັນ
Trans-1,2-Dichloroethene	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
1.1-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Butanone	11 UJ	11 U	11 U	11 U	11 U	11 U	11 U	11 UJ
Cis-1,2-Dichloroethene	4 U	4 U	4 U	4 U	4 U	4 U	4 U	4 U
Methyl tert-Butyl Ether	1 UJ	1 U	1 U	1 U	1 U	1 U	1 U	1 UJ
1,2-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chioroform	3 U	3 U	3 U	3 U	3 U	3 U	3 U	3 U
Carbon Tetrachioride	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
1,1,1-Trichloroethane	2 U	2 U	2 U	2 U	1 J	2 J	2 U	0.9 J
Benzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichleropropane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromodichioromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Cis-1,3-Dichioropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trans-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1,2-Trichloroethane	1 U	1 U	1 U	1 U	1 U	1 0	1 U	1 U
Dibromochloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	1 U	1 U	1 U	1 U	1 U	1 0	1 U	1 U
4-Methyl-2-pentanone	5 U	5 U 0.8 J		5 U	5 U	5 U	5 U	5 U
Toluene Tetrachloroethene	1 0	0.8 J	1 J 1 U	1 U	0.6 J 1 U	0.8 J 1 U	1 J	2 J 1 U
2-Hexanone	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U
1.2-Dibromoethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chlorobenzene	1 U	1 U	1 U	1 U	1 Ü	1 U	1 U	1 U
Ethylbenzene Ethylbenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 0
m-2/or p-Xylene	0.6 J	1 U	1 0	1 0	1 0	1 U	1 U	1 U
o-xylene	1 U	1 U	1 0	1 0	1 0	1 U	1 U	1 0
Xylenes (total)	1			<del>'</del>				
Styrene	3 UJ	3 U	3 U	3 U	3 U	3 U	3 U	3 UJ
1,1,2,2-Tetrachloroethane	1 U	1 U	1 U	1 0	1 U	1 0	1 U	1 U
1,3-Dichiorobenzene	1 U	1 U	1 U	1 0	1 U	1 U	1 U	1 0
1.A-Dichlorobenzene	1 U	1 U	1 U	10	1 0	1 0	1 0	10
1,2-Dichlorobenzene	1 0	1 0	1 U	1 0	1 0	1 U	1 0	10
1,2,4-Trichlorobenzene	<del></del>	<del></del>	<del></del> _	<del></del>				
1,2-Dibromo-3-chloropropane			<del>                                     </del>					
Bromochloromethane								
Francisco de la constancia de la constan	L							

Sample ID:	CPT03-06	CPT-03-07	CPT-04-01	CPT-04-02	CPT-04-02DUP	CPT-04-02	CPT04-02DP	CPT-04-03
Sample Number:	EABX5					EABW5	EABW6	
Sample Pate:	6/2/00	6/2/00	5/29/00	5/29/00	5/29/00	5/29/00	5/29/00	5/29/00
Laboratory:	MitKem	ESAT	ESAT	ESAT	ESAT	MitKem	MitKem	ESAT
<u> </u>	92	102	32	46	46	46	46	56
Sample Screen Depth (ft below ground): Units:	ug/L	ug/L	wg/L	ug/L	ug/L	ug/L	ug/L	ug/L
	uyrc	- Opt	- upc	ugic		- Ugra	- dy.L	y
Parameter:		1 U	1 U	1 U	1 U			1 U
Dichlorodifluoromethane chloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
	1 0	1 U	1 U	1 U	1 0	1 0	1 0	1 U
Vinyl Chloride  Bromomethane	1 U	1 U	1 0	1 0	1 0	1 U	1 U	1 0
Chloroethane	1 U	1 U	1 U	1 U	1 0	1 U	1 U	1 U
Trichlorofluoromethane	<del></del>	1 U	1 U	1 U	1 0			1 U
1,1,2-Trichloro-1,2,2-trifluoromethane	<del> </del>	2 UJ	2 UJ	2 UJ	2 UJ			2 UJ
1,1-Dichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 Ü	1 U
Acetone	5 U	2 UJ	2 U	10 B	2 U	5 U	5 U	2 U
Carbon Disulfide	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Methylene Chioride	2 U	1 U	1 U	1 U	1 U	2 U	2 U	1 U
Methyl Acetate		1 UJ	1 UJ	1 UJ	1 UJ	_	_	1 UJ
Trans-1,2-Dichloroethene	1 U	2 U	2 U	2 U	2 U	1 U	1 U	2 U
1,1-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Butanone	5 U	11 UJ	11 U	11 U	11 U	5 U	5 U	11 U
Cis-1,2-Dichloroethene	1 U	4 U	4 U	4 U	4 U	1 U	1 U	4 U
Methyl tert-Butyl Ether		1 UJ	1 U	1 U	1 U			1 U
1,2-Dichloroethane	1 U	1 Ü	1 U	1 U	1 U	1 U	1 U	1 U
Chloroform	1 U	3 U	3 U	3 U	3 U	1 U	1 U	3 U
Carbon Tetrachloride	1 U	2 U	2 U	2 U	2 U	1 U	1 U	2 U
1,1,1-Trichloroethane	0.9 J	0.8 J	2 U	2 U	2 U	1 U	1 U	2 U
Senzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichloropropane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromodichloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Cis-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trans-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1,2-Trichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Dibromochloromethane	1 U	1 U	1 U	1 U	1 U	<u>1 U</u>	1 U	1 U
Bromoform	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 0
4-Methyl-2-pentanone	5 U	5 U 2 J	5 U	5 U 0.8 J	5 U	5 U	5 U	5 U
Toluene	1 U	1 U	1 J 1 U	1 U	0.9 J	0.6 J 1 U	0.8 J 1 U	0.8 J 1 U
Tetrachloroethene 2-Hexanone	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U
1.2-Dibromoethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 0
Chlorobenzene	1 0	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Ethylbenzene	1 U	1 U	1 0	1 U	1 U	1 U	1 U	1 U
m-&/or p-Xylene		1 U	1 0	1 0	1 0			1 0
o-xylene		1 U	1 U	1 U	1 0			1 0
Xylenes (total)	1 U					1 U	1 U	
Styrene	1 U	3 ÜJ	3 U	3 U	3 U	1 U	1 U	3 U
1,1,2,2-Tetrachloroethane	1 0	1 U	1 U	1 Ü	1 0	1 U	1 Ü	1 U
1,3-Dichlorobenzene	1 Ü	1 U	1 U	1 0	1 U	1 U	1 U	1 U
1,4-Dichlorobenzene	1 U	1 U	1 U	1 0	1 0	1 U	1 0	1 U
1,2-Dichlorobenzene	1 U	1 U	1 Ü	1 Ü	1 U	1 U	1 0	1 U
1,2,4-Trichlorobenzene	1 U					1 U	1 U	
1,2-Dibromo-3-chloropropane	1 R	-				1 R	1 R	
Bromochioromethane	1 U			<u> </u>	<del> </del>	1 U	10	
Control of the contro					<del></del>	·		

Section: 5 Revision: 1 Date: 28 March 2001 Page: 14 of 35

Sample ID:	CPT-04-04	CPT-04-05	CPT-04-06	CPT-04-07	CPT-05-01	CPT-05-02	CPT-05-03	CPT-05-04
Sample Number:					_			_
Sample Date:	5/29/90	5/29/00	5/29/00	5/29/00	5/29/00	5/29/00	5/29/00	5/29/00
Laboratory:	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT
Sample Screen Depth (ft below ground):	71	78	84	93	35	43	51	57
Units:	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Parameter:				·			<u> </u>	
Dichlorodifluoromethane	1 U		1 U	1 0	1 U	1 U	1 U	1 U
chloromethane	1 U		1 U	1 U	1 U	1 U	1 U	1 U
Vinyi Chloride	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromomethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichlorofluoromethane	1 U	1 U	1 U	1 U	1 U	1 Ú	1 U	1 U
1.1.2-Trichloro-1.2.2-triffuoromethane	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ
1.1-Dichloroethene	1 U	1 U	1 U	1 U	1 U	1 Ú	1 U	1 U
Acetone	2 U	2 U	8 B	9 B	2 U	2 U	2 U	2 U
Carbon Disulfide	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Methylene Chioride	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Methyl Acetate	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ
Trans-1,2-Dichloroethene	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
1,1-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Butanone	11 U	11 Ü	11 U					
Cis-1,2-Dichloroethene	4 U	4 U	4 U	4 U	4 U	4 U	4 U	4 U
Methyl tert-Butyl Ether	1 U	1 U	1 U	1 U	J 1	1 U	1 U	1 U
1,2-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroform	3 U	3 U	3 U	3 U	3 U	3 U	3 U	3 U
Carbon Tetrachioride	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
1,1,1-Trichioroethane	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
Benzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichioroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichloropropane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromodichioromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Cis-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trans-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 Ü	1 U_	1 U	1 U
1,1,2-Trichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Dibromochioromethane	1 U	1 U	1 U	1 U	1 Ü	1 U	1 U	1 U
Bromoform	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
4-Methyl-2-pentanone	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U
Toluene	0.7 J	1 J	0.8 J	1 J	1 J	1 J	1 J	1 J
Tetrachioroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Hexanone	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U
1,2-Dibromoethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Ethylbenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-8/or p-Xylene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
o-xylene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Xylenes (total)	3 U	3 U	3 U	3 U	3 U	3 11	3 11	<del></del> _
Styrene 1,1,2,2-Tetrachioroethane	1 U	1 U	1 U	1 U		3 U	3 U	3 U
1.3-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1.4-Dichlorobenzene	1 0	1 U	1 U	1 U	1 0		1 U	1 U
	1 U	1 0	1 U	1 U		1 U	1 υ	1 U
1,2-Dichlorobenzene		†			1 U	1 U	1 U	1 U
1,2,4-Trichlorobenzene	ļ <del></del>	<del></del>						
1,2-Dibromo-3-chloropropane								
Bromochloromethane						***		

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Sample ID:	CPT-05-05	CPT-05-05DUP	CPT-05-06	CPT-05-07	CPT-06-01	CPT-06-02	CPT-06-03	CPT-06-04
Sample Number:	<b> </b>							
Sample Date:	5/30/00	5/30/00	5/30/00	5/30/00	6/1/00	6/1/00	6/1/00	6/1/00
Laboratory:	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT
Sample Screen Depth (ft below ground):	69.5	69.5	78	87	35	42	63	62
Units:	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Parameter:								
Dichlorodifluoromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
chloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Vinyl Chloride	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromomethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichlorofluoromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1,2-Trichloro-1,2,2-triffuoromethane	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ
1,1-Dichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Acetone	2 U	2 U	100 J	2 U	2 U	53 J	11 J	7 J
Carbon Disulfide	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Methylene Chioride	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Methyl Acetate	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ
Trans-1,2-Dichloroethene	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
1,1-Dichloroethane	1 U	1 U	1 U 16 J	1 U	1 U	1 U	1 U	1 U
2-Butanone	<del></del>	·	4 U	4 U	4 U	4 U		
Cis-1,2-Dichloroethene	4 U	4 U					4 U	4 U
Methyl tert-Butyl Ether	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichloroethane	3 U	3 U	3 U	3 U	1 U 3 U	3 U	3 U	1 U
Chloroform								3 U
Carbon Tetrachloride	2 U	2 U	2 U 2 U	2 U	2 U	2 U 2 U	2 U	2 U 2 U
1,1,1-Trichloroethane	1 U	1 U	1 U	1 0	1 U	1 U	1 U	1 U
Benzene Trichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
	1 0	1 U	1 U	1 0	1 U	1 U	1 U	1 0
1,2-Dichloropropane Bromodichloromethane	1 1 0	1 U	1 U	1 U	1 0	1 U	1 U	1 0
Cis-1,3-Dichloropropene	1 0	1 U	1 0	1 0	1 U	1 U	1 U	1 U
Trans-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 0
1.1.2-Trichloroethane	1 0	1 0	1 U	1 U	1 0	1 U	1 U	10
Dibromochloromethane	1 0	1 0	1 U	1 0	1 0	1 0	1 U	1 0
Bromoform	1 0	1 0	1 U	1 0	1 0	1 U	1 U	1 0
4-Methyl-2-pentanone	5 U	5 U	5 U	5 Ü	5 U	5 U	5 U	5 U
Toluene	1 1	0.6 J	0.8 J	1 J	0.6 J	0.8 J	0.8 J	1 1
Tetrachloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 Ü
2-Hexanone	5 U	5 U	5 U	5 U	5 U	5 U	5 Ü	5 U
1,2-Dibromoethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chlorobenzene	1 U	1 0	1 U	1 U	1 U	1 U	1 U	1 U
Ethylbenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-&/or p-Xylene	1 U	1 U	1 U	1 U	0.7 J	1 U	1 U	1 U
o-xylene	1 U	1 U	1 U	1 U	1 U	1 U	1 Ü	1 U
Xylenes (total)	-							
Styrene	3 U	3 U	3 U	3 U	3 UJ	3 UJ	3 UJ	3 UJ
1,1,2,2-Tetrachioroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,3-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 Ü
1,4-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 Ü	1 0
1,2,4-Trichlorobenzene								
1,2-Dibromo-3-chloropropane			-					
Bromochloromethane								
<u> </u>	·				<u> </u>	<del> </del>		

Table 5-1
Groundwater Analytical Results - CPT - VOCs
Evergreen Manor, Roscoe, Illinois

Sample ID:	CPT-06-05	CPT-06-06	CPT06-06	CPT-06-07	CPT-09-01	CPT-09-02	CPT-09-03	CPT-09-04
Sample Number:			EABW9			-	_	-
Sample Date:	6/1/00	6/1/00	6/1/00	6/1/00	5/30/00	5/30/00	5/30/00	5/30/00
Laboratory:	ESAT	ESAT	MitKem	ESAT	ESAT	ESAT	ESAT	ESAT
Sample Screen Depth (ft below ground):	74	85	85	92	35	45	55	68
Units:	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Parameter:	<u></u>		<del></del>				·	
Dichlorodifluoromethane	1 U	1 U		1 U	1 U	1 U	1 U	1 U
chloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	J 1
Vinyt Chloride	1 U	1 U	1 U	1 U	1 Ü	1 U	1 U	1 U
Bromomethane	1 U	1 U	1 U	1 Ü	1 U	1 U	1 U	1 U
Chloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	י נ
Trichlorofluoromethane	1 υ	1 υ		1 υ	1 U	1 υ	1 υ	1 υ
1,1,2-Trichloro-1,2,2-trifluoromethane	2 UJ	2 UJ		2 UJ	2 UJ	2 UJ	2 UJ	2 UJ
1,1-Dichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Acetone	11 J	2 U	7 U	5 J_	2 U_	2 U	2 U	2 U
Carbon Disulfide	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Methylene Chloride	1 U	1 U	2 U	1 U	1 U	1 U	1 U	1 U
Methyl Acetate	1 UJ	1 UJ		1 UJ	1 UJ	1 UJ	1 UJ	1 UJ
Trans-1,2-Dichloroethene	2 U	2 U	1 U	2 U	2 U	2 U	2 U	2 U
1,1-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Butanone	11 U	11 U	5 U	11 U	11 U	11 U	11 U	11 U
Cis-1,2-Dichloroethene	4 U	4 U	1 U	4 U	4 U	4 U	4 U	4 U
Methyl tert-Butyl Ether	1 U	1 U		1 U	1 U	1 U	1 U	1 U
1,2-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroform	3 U	3 U	1 Ü	3 U	3 U	3 U	3 U	3 U
Carbon Tetrachloride	2 U	2 U	1 U	2 U	2 U	2 U	2 U	2 U
1,1,1-Trichloroethane	2 U	2 U	1 0	2 U	2 U	2 U	2 U	2 U
Benzene	1 U	1 U	1 U	1 U	0.5 J	1 U	1 U	1 U
Trichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichloropropane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromodichloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Cis-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trans-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1,2-Trichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 0
Dipromochloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	1 U	1 U	1 U 5 U	1 U	1 U	1 U	1 U	1 U
4-Methyl-2-pentanone	1 J	5 U		5 U	5 U	5 U	5 U	5 U
Toluene Tetrachioroethene	1 0	2 J 1 U	1 U	1 J	1 J	0.8 J 1 U	0.8 J 1 U	0.8 J 1 U
2-Hexanone	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U
1,2-Dibromoethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chlorobenzene	1 U	1 U	1 U	1 0	1 U	1 U	1 U	1 U
Ethylbenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-8/or p-Xylene	1 U	1 0	<del></del> -	1 U	1 0	1 U	1 U	1 U
o-xylene	1 0	1 0		1 0	1 U	1 U	1 0	1 U
Xylenes (total)	<del> </del>		1 U	- '-				
Styrene	3 UJ	3 UJ	1 U	3 UJ	3 U	3 U	3 U	3 U
1,1,2,2-Tetrachloroethane	1 0	1 U	1 0	1 U	1 U	1 U	1 U	1 U
1,3-Dichlorobenzene	1 0	1 0	1 0	1 0	1 0	1 U	1 0	1 U
1,4-Dichlorobenzene	1 0	1 U	1 0	1 0	1 0	1 U	1 U	1 U
1,2-Dichiorobenzene	1 0	1 U	1 U	1 0	1 U	1 0	1 U	1 U
1,2,4-Trichlorobenzene	<del>                                     </del>		1 0					
1,2-Dibromo-3-chloropropane	<del>                                     </del>	<del></del>	1 R		<del></del>			
Bromochioromethane	+		1 0		<del>                                     </del>			
Camaramana								

Parala ID.	CPT-09-05	CPT-09-06	CPT-09-07	CPT-10-01	CPT-10-02	CPT-10-03	CPT-10-04	CPT-10-05
Sample ID:	CF1-03-03	CF1-03-06	CF1-03-07	CF 1-10-01		<u> </u>		
Sample Number:		57000	50000	5/31/00	5/31/00	5/31/00	6/1/00	€/1/00
Sample Date:	5/30/00	5/30/00	5/30/00	ESAT	ESAT	ESAT	ESAT	ESAT
Laboratory:	ESAT	ESAT	ESAT 88	25	36	42	55	65
Sample Screen Depth (ft below ground):	75	85						
Units:	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Parameter:							4 11	4 11
Dichlorodifluoromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
chloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Vinyl Chloride	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromomethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chioroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichlorofluoromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1,2-Trichloro-1,2,2-trifluoromethane	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ
1,1-Dichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Acetone	2 U	11 J	8 J	2 U	2 U	40 B	2 U	2 U
Carbon Disulfide	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Methylene Chloride	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Methyl Acetate	1 UJ	1 UJ	. 1 UJ	1 U	1 U	1 U	1 UJ	1 Ü
Trans-1,2-Dichloroethene	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
1,1-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Butanone	11 U	11 U	11 U	11 U	11 U	11 U	11 U	11 U
Cis-1,2-Dichloroethene	4 U	4 U	4 U	4 U	4 U	4 U	4 U	4 U
Methyl tert-Butyl Ether	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1.2-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroform	3 U	3 U	3 U	3 U	3 U	3 U	3 U	3 U
Carbon Tetrachloride	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
1,1,1-Trichloroethane	2 U	2 U	2 U	2 U	1 J	2 U	1 J	0.8 J
Benzene	1 Ü	1 U	0.6 J	1 U	1 U	1 U	1 U	1 U
Trichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichloropropane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromodichloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Cis-1,3-Dichloropropene	1 U	1 Ü	1 U	1 U	1 U	1 U	1 U	1 U
Trans-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1,2-Trichloroethane	1 U	1 U	1 U	1 U	1 Ú	1 U	1 U	1 U
Dibromochloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
4-Methyl-2-pentanone	5 U	5 U	5 U	5 Ü	5 U	5 U	5 U	5 U
Toluene	0.5 J	0.8 J	2 J	0.9 J	1 U	1 J	0.5 J	0.9 J
Tetrachioroethene	1 U	1 U	1 U	1 U	0.9 J	1 U	0.6 J	1 U
2-Hexanone	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U
1,2-Dibromoethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Ethylbenzene	1 U	1 U	0.6 J	1 U	1 U	1 U	1 U	1 U
m-&/or p-Xylene	1 U	1 U	0.7 J	1 U	1 U	1 U	1 Ü	1 U
o-xylene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Xylenes (total)								
Styrene	3 U	3 U	3 U	3 U	3 U	3 U	3 UJ	3 U
1,1,2,2-Tetrachloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,3-Dichlorobenzene	1 U	1 U	1 0	1 U	1 U	1 U	1 U	1 U
1,4-Dichlorobenzene	1 0	1 U	1 0	1 U	1 U	1 U	1 U	1 U
1,2-Dichlorobenzene	1 0	1 0	1 0	1 U	1 U	1 U	1 U	1 U
1,2,4-Trichlorobenzene				<del></del>				
1,2-Dibromo-3-chloropropane								
Bromochloromethane	<del></del>					<del></del>		
	<u> </u>			1				

Sample ID:	CPT-10-06	CPT-10-07	CPT-10-4DUP	CPT-11-01	CPT11-01	CPT-11-02	CPT-11-03	CPT-11-04
Sample Number:		-			EABX7		_	1
Sample Date:	6/1/00	6/1/00	6/1/00	6/3/00	6/3/00	6/3/00	6/3/00	6/3/00
Laboratory:	ESAT	ESAT	ESAT	ESAT	MitKem	ESAT	ESAT	ESAT
Sample Screen Depth (ft below ground):	73	90	55	45	45	58	70	81
Units:	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Parameter:			·	<del></del>	<u> </u>		-	
Dichlorodifluoromethane	1 U	1 U	1 U	1 U		1 U	1 U	1 U
chloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Vinyl Chloride	1 U	1 U	1 U	1 UJ	1 U	1 UJ	1 UJ	1 UJ
Bromomethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichlorofluoromethane	1 U	1 U	1 U	1 υ	_	1 U	1 U	1 U
1.1.2-Trichioro-1.2.2-trifluoromethane	2 UJ	2 UJ	2 UJ	2 UJ		2 UJ	2 UJ	2 UJ
1.1-Dichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Acetone	24 J	2 U	2 U	2 UJ	5 U	2 UJ	9 J	2 UJ
Carbon Disulfide	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Methylene Chloride	1 U	1 U	1 U	1 U	2 U	1 U	1 U	1 U
Methyl Acetate	1 UJ	1 UJ	1 UJ	1 UJ	_	1 UJ	1 UJ	1 00
Trans-1,2-Dichloroethene	2 U	2 U	2 U	2 U	1 U	2 U	2 U	2 U
1,1-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Butanone	11 U	11 U	11 U	11 UJ	5 U	11 UJ	11 UJ	11 UJ
Cis-1,2-Dichloroethene	4 U	4 U	4 Ü	4 U	1 U	4 U	4 U	4 U
Methyl tert-Butyl Ether	1 U	1 U	1 U	1 UJ		1 UJ	1 UJ	1 UJ
1,2-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroform	3 U	3 U	3 U	3 U	1 U	3 U	3 U	3 U
Carbon Tetrachioride	2 U	2 U	2 U	2 U	1 U	2 U	2 U	2 Ü
1,1,1-Trichloroethane	2 U	2 U	1 J	2 U	1 U	2 U	0.8 J	0.8 J
Benzene	1 U	1 U	1 U	1 U	1 Ü	1 U	1 U	1 U
Trichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichloropropane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromodichloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U .
Cis-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trans-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1,2-Trichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Dibromochloromethane	1 U	1 U	1 U	1 U	U 1	1 U	1 U	1 U
Bromoform	1 0	1 U	1 U	1 UJ	1 U	1 UJ	1 UJ	1 UJ
4-Methyl-2-pentanone	5 U	5 U	5 U	5 UJ	5 U	5 UJ	5 UJ	5 UJ
Toluene	1 J	1 J	1 U	1 U	1 U	1 J	1 J	0.9 J
Tetrachioroethene	1 U	1 U	0.6 J	1 U	1 U	1 U	1 U	1 U
2-Hexanone	5 U	5 U	5 U	5 UJ	5 U	5 UJ	5 UJ	5 UJ
1,2-Dibromoethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Ethylbenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-8/or p-Xylene	1 U	1 U	1 U	1 U		1 U	1 U	1 U
o-xylene	1 U	1 U	1 U	1 U		1 U	1 U	1 U
Xylenes (total)					0.6 J			
Styrene	3 UJ	3 UJ	3 UJ	3 U	1 U	3 U	3 U	3 U
1,1,2,2-Tetrachloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,3-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 Ü	1 U
1,4-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2,4-Trichlorobenzene		<del>-</del>			1 U			
1,2-Dibromo-3-chloropropane					1 R			
Bromochioromethane				***	1 U			

Page. 19 of 35

Sample Number:	S	CPT-11-05	CPT11-05	CPT11-05DP	CPT-11-06	CPT-11-07	CPT11-08	CPT-11-08	CPT-12-01
Sample Date:	<del></del>	CF1-11-05			CPI-II-O	CFI-11-07		CF1-11-06	CF1-12-01
Laboratory:   SAAT   MRIStone   SAAT   MRIStone   SAAT   SAAT		come			6000	67700		£7200	6/5/00
Sample Screen Depth (it below ground):   93   93   93   93   92   114   125   125   45									
Unite:									
Dichlorodifuctomethane		<del></del>							
Dehlorodiflucromethane		ugrL	ug/L	eg/L	ug/L	ug/L	ug/L	ugit	ugvi
New Comments		1 11			4 11	T 4 11		4 11	
Very   Chlorids							1 11		
Stromomethane									
Shiorsethane									
Tickhloroffuoromethane									
1,1,2-Trichloro-1,2,2-driffuoromethane			1 0	1 0					
1.1-Dichloroethene		<u> </u>		<del></del>					
Acetons							4 11		
Sarbon Disulfide									
Methylane Chloride		<del></del>							
Methyl Acetate									
Trans-1,2-Dichloroethene			20						
1.1-Dichloroethane			1 11				1 11		
2-Butanone									
Cis-12-Dichloroethene									
Methyl tan-Butyl Ether									
1		1				<u> </u>			
Chloroform				L			<u> </u>		
Carbon Tetrachloride							<u> </u>		
1,1,1-Trichloroethane									
Benzene									
Trickloroethene	<del></del>								
1.2-Olchloropropane								<del></del>	
Stromodichloromethane	<del></del>								
Cis-1,3-Dichloropropene									
Trans-1,3-Dichloropropene									
1,1,2-Trichloroethane									
Dibromochloromethane									
Bromoform	<del></del>								
##Methyl-2-pentanone		1 UJ							
Toluene		5 UJ				5 UJ			
2-Hexanone		1 J	2		1 J	2 J			
1,2-Olbromoethane         1 U	Tetrachloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chlorobenzene	2-Hexanone	5 UJ	5 U	5 U	5 UJ	5 UJ	5 U	5 UJ	5 U
Ethylbenzene	1,2-Dibromoethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-∨ p-Xylene         1 U           1 U         1 U          1 U         1 U          1 U	Chlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
o-xylene         1 U          1 U         1 U          1 U         1	Ethylbenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Xylenes (total)	m-&/or p-Xylene	1 U			1 U	1 U		1 U	1 U
Styrene	o-xylene	1 U	***		1 U	1 U	_	1 U	1 U
1,1,2,2-Tetrachloroethane         1 U <th>Xylenes (total)</th> <th></th> <th>1 U</th> <th>1 U</th> <th></th> <th></th> <th>1 U</th> <th></th> <th></th>	Xylenes (total)		1 U	1 U			1 U		
1,3-Dichlorobenzene         1 U	Styrene		1 U	1 U	3 U	3 U	1 U	3 U	3 U
1,4-Dichlorobenzene         1 U	1,1,2,2-Tetrachloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichlorobenzene         1 U	1,3-Dichlorobenzene		1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2,4-Trichlorobenzene — 1 U 1 U — 1 U — 1 U 1,2-Dibromo-3-chloropropane — 1 R 1 R 1 R 1 R 1 R	1,4-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dibromo-3-chloropropane 1 R 1 R 1 R 1 R	1,2-Dichlorobenzene	1 U		1 U	1 U	1 U	1.0	1 U	1 U
	1,2,4-Trichlorobenzene								
Bromochloromethane 1 U 1 U 1 U 1 U	1,2-Dibromo-3-chloropropane						1 R	_	
	Bromochloromethane		1 U	1 U		_	1 U	_	

Sample ID:	CPT-12-01DUP	CPT-12-02	CPT-12-03	CPT-12-03	CPT-12-04	CPT-12-04BUP	CPT-12-05	CPT12-06
Sample Number:				EABQ3				EABQ7
Sample Date:	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00
Laboratory:	ESAT	ESAT	ESAT	MitKem	ESAT	ESAT	ESAT	MitKem
Sample Screen Depth (ft below ground):		62	70	70	81	81	93	102
Units:	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Parameter:								
Dichlorodifluoromethane	1 U	1 U	1 U		1 U	1 Ú	1 U	
chloromethane	1 U	1 U	1 Ü	1 U	1 U	1 Ü	1 Ü	1 U
Vinyl Chloride	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromomethane	1 0	1 Ü	1 U	1 U	1 U	1 U	1 U	1 U
Chloroethane	1 0	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichlorofluoromethane	1 U	1 U	1 U	_	1 Ü	1 U	1 U	
1.1.2-Trichloro-1,2,2-trifluoromethane	2 U	2 U	2 U		2 U	2 U	2 U	
1,1-Dichloroethene	1 U	1 Ü	1 U	1 U	1 U	1 U	1 U	1 U
Acetone	1 B	1 B	2 B	5 U	2 <b>B</b>	7 B	2	5 U
Carbon Disulfide	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Methylene Chloride	1 U	1 U	1 U	2 U	1 U	1 U	1 U	2 U
Methyl Acetate	1 UJ	1 UJ	1 UJ		1 UJ	1 UJ	1 UJ	_
Trans-1,2-Dichloroethene	2 U	2 U	2 U	1 U	2 U	2 U	2 U	1 U
1,1-Dichloroethane	1 U	1 U	1 Ú	1 U	1 Ú	1 U	1 U	1 U
2-Butanone	11 U	11 U	11 U	5 U	11 U	11 U	11 U	5 U
Cis-1,2-Dichloroethene	4 U	4 U	4 U	1 U	4 Ü	4 U	4 U	1 U
Methyl tert-Butyl Ether	1 U	1 U	1 U		1 U	1 U	1 U	
1,2-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroform	3 U	3 U	3 U	1 U	3 U	3 U	3 U	1 U
Carbon Tetrachioride	2 U	2 U	2 U	1 U	2 U	2 U	2 U	1 U
1,1,1-Trichioroethane	2 U	2 U	2 U	1 U	2 U	2 U	2 U	1 U
Benzene	1 U	1 U	1 U	1 U	1 U	0.5 J	1 U	1 U
Trichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichloropropane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromodichioromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Cis-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 Ü	1 U	1 U	1 Ü
Trans-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1,2-Trichloroethane	1 0	1 U	1 0	1 U	1 U	1 U	1 U	1 U
Dibromochloromethane	1 0	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	5 U	5 U	5 U	5 U	5 Ü	5 U	5 U	5 U
4-Methyl-2-pentanone Toluene	0.5 J	1 U	0.9 J	0.8 J	0.9 J	1 0	2 J	1
Tetrachloroethene	1 0	1 U	1 U	1 U	1 U	1 0	1 U	1 U
2-Hexanone	5 U	5 U	5 U	5 Ü	5 U	5 U	5 U	5 U
1.2-Dibromoethane	1 0	1 U	1 U	1 0	1 U	1 0	1 U	1 U
Chlorobenzene	1 0	1 U	1 U	1 Ŭ	1 0	1 0	1 0	1 U
Ethylbenzene	1 0	1 U	1 U	1 Ü	1 0	1 0	1 0	1 U
m-&/or p-Xylene	1 0	1 U	1 Ü		1 U	1 0	1 Ŭ	
o-xylene	1 U	1 U	1 U		1 Ü	1 Ü	1 0	_
Xylenes (total)				1 U				1 U
Styrene	3 U	3 U	3 U	1 U	3 U	3 U	3 U	1 0
1,1,2,2-Tetrachloroethane	1 U	1 U	1 Ü	1 U	1 Ü	1 0	1 U	1 U
1,3-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,4-Dichlorobenzene	1 U	1 U	1 U	1 U	1 Ü	1 0	1 0	1 U
1,2-Dichlorobenzene	1 U	1 U	1 Ü	1 Ü	1 Ü	1 U	1 0	1 Ü
1,2,4-Trichlorobenzene				1 Ü				1 0
1,2-Dibromo-3-chloropropane				1 R				1 R
Bromochloromethane				1 Ü				1 0
						<u> </u>		

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Sample ID:	CPT-12-06	CPT-12-07
Sample Number:		
Sample Date:	6/5/00	6/5/00
Laboratory:	ESAT	ESAT
Sample Screen Depth (ft below ground):	102	118
Units:	ug/L	ug/L
Parameter:	UyrE	
	1 U	1 U
Dichlorodifluoromethane	1 U	1 U
chloromethane	1 U	1 U
Vinyl Chloride	1 1 0	
Bromomethane		
Chioroethane	1 U	1 U
Trichiorofluoromethane	1 U	1 U
1,1,2-Trichloro-1,2,2-triffuoromethane	2 U	2 U
1,1-Dichloroethene	1 U	1 U
Acetone	2	2
Carbon Disulfide	1 U	1 U
Methylene Chioride	1 U	1 U
Methyl Acetate	1 UJ	1 UJ
Trans-1,2-Dichloroethene	2 U	2 U
1,1-Dichloroethane	1 U	1 U
2-Butanone	11 U	11 U
Cis-1,2-Dichloroethene	4 U	4 U
Methyl tert-Butyl Ether	1 U	1 U
1.2-Dichloroethane	1 U	1 U
Chloroform	3 U	3 U
Carbon Tetrachloride	2 U	2 Ü
1,1,1-Trichloroethane	2 U	2 U
Benzene	1 0	1 0
Trichloroethene	1 U	1 0
1,2-Dichloropropane	1 0	1 0
Bromodichioromethane	1 0	1 0
Cis-1,3-Dichloropropene	1 0	1 0
	1 0	1 0
Trans-1,3-Dichloropropene	1 0	1 U
1,1,2-Trichioroethane		
Dibromochloromethane	1 U	1 U
Bromoform	1 U	1 0
4-Methyl-2-pentanone	5 0	5 U
Toluene	2 J	2 J
Tetrachioroethene	1 U	1 U
2-Hexanone	5 U	5 U
1,2-Dibromoethane	1 U	1 U
Chlorobenzene	1 U	1 U
Ethylbenzene	1 U	1 U
m-&/or p-Xylene	1 U	1 U
o-xylene	1 U	1 U
Xylenes (total)		
Styrene	3 U	3 U
1,1,2,2-Tetrachioroethane	1 U	1 U
1,3-Dichlorobenzene	1 U	1 U
1,4-Dichlorobenzene	1 U	1 U
1,2-Dichlorobenzene	1 U	1 U
1,2,4-Trichlorobenzene		
1,2-Dibromo-3-chloropropane		
	1	+

Note:

- U Constituent not detected; method detection limit (MDL) of the analysis reported.
- J Concentration reported is an estimated value.
- UJ The analyte was not detected above the reported sample quantition limit.
- B Designates the constituent was detected in the method blank.
- R The sample results are rejected due to deficiencies in the ability to analyze the sample and
- --- Indicates compound not analyzed.

#### Table 5-2 Groundwater Analytical Results - Monitoring Wells - VOCs Evergreen Manor, Roscoe, Illinois

Sample ID:	MW101D	MW1018	MW102D	MW1025	MW103D	MW103D-DL	MW1035	MW104D
Sample Number:		_	_	_	_	_		_
Sample Date:	8/30/00	6/30/00	8/1/00	8/1/00	5/31/00	5/31/00	5/31/00	8/2/00
Laboratory:	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT
Units	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
PARAMETER	<u> </u>					<del></del>		
1,1,1-Trichloroethane	2 U	2 U	2 U	2 U	3 J	50 U	2 U	1 J
1,1,2,2-Tetrachioroethane	1 U	1 U	1 U	1 U	1 U	25 U	1 U	1 U
1.1.2-Trichloro-1,2,2-trifiuoromethane	2 UJ	2 UJ	2 U	2 UJ	300 J	180 J	2 J	2 UJ
1,1,2-Trichioroethane	1 U	1 U	1 υ	1 υ	1 υ	25 U	1	1 U
1,1-Dichloroethane	1 U	1 U	1 U	1 U	1 U	25 U	1 U	1 U
1.1-Dichioroethene	1 U	1 U	1 U	1 U	1 υ	25 U	1 U	1 U
1.2.4-Trichlorobenzene		_	_	_				-
1,2-Dibromo-3-chioropropane	_							-
1,2-Dibromoethane	1 U	1 U	1 U	1 U	1 U	25 U	1 Ü	1 U
1,2-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	25 U	1 U	1 U
1,2-Dichloroethane	1 U	1 U	1 U	1 U	1 U	25 U	1 Ü	1 U
1,2-Dichloropropane	1 U	1 U	1 U	1 U	1 U	25 U	1 U	1 υ
1,3-Dichlorobenzene	1 Ú	1 U	1 U	1 U	1 U	25 U	1 U	1 U
1,4-Dichiorobenzene	1 U	1 U	1 U	1 U	1 U	25 U	1 U	1 U
2-Butanone	11 U	11 U	11 U	11 U	11 U	250 U	11 U	11 UJ
2-Hexanone	5 Ü	5 Ū	5 U	5 U	5 U	125 U	5 U	5 U
4-Methyl-2-pentanone	5 U	5 U	5 U	5 U	5 U	125 U	5 U	5 U
Acetone	2 U	2 U	2 U	2 U	35 B	50 U	2 U	2 UJ
Benzene	1 U	1 Ü	1 U	1 U	1 U	25 U	1 U	1 U
Bromochioromethane							-	
Bromodichioromethane	1 U	1 U	1 U	1 U	ŤÜ	25 U	1 U	1 U
Bromoform	1 U	1 U	1 U	1 U	1 U	25 U	1 U	1 U
Bromomethane	1 U	1 U	1 U	1 U	1 U	25 U	1 U_	1 U
Carbon Disuifide	1 U	1 U	1 U	1 U	1 U	25 U	1 U	1 U
Carbon Tetrachioride	2 U	2 U	2 U	2 U	2 U	50 U	2 U	2 U
Chlorobenzene	1 U	1 U	1 U	1 U	1 Ü	25 U	1 U	1 U
Chloroethane	1 U	1 U	1 U	1 U	1 U	25 U	1 U	1 U
Chloroform	3 U	3 U	3 U	3 U	3 U	75 U	3 υ	3 U
Chloromethane	1 U	1 U	1 U	1 U	1 U	25 U	1 U	1 U
Cls-1,2-Dichloroethene	4 U	4 U	4 U	4 U	4 U	100 U	4 U	4 U
Cls-1,3-Dichioropropene	1 U	1 U	1 U	1 U	1 U	25 U	1 U	1 U
Dibromochioromethane	1 U	1 U	1 U	1 U	1 U	25 U	1 U	1 U
Dichlorodifluoromethane			1 U			25 U	1 U	1 U
Ethylbenzene	1 U	1 U	1 0	1 U	1 U	25 U	1 U	1 U
m-8/or p-Xylene	1 1111	1 UJ	1 0	1 W	1 0	25 U	1 U	
Methyl Acetate	1 U	1 U	1 U	1 U	1 U	25 U	1 UJ 1 U	1 UJ 1 UJ
Methyl tert-Butyl Ether	1 0	1 0	1 U	1 0	1 0	25 U	1 U	1 00
Methylene Chioride o-xylene	10	1 0	10	1 0	1 0	25 U	1 U	10
	3 U	3 U	3 U	3 UJ	3 UJ	75 UJ	3 UJ	3 UJ
Styrene Tetrachioroethene	1 0	1 0	1 U	1 U	0.5 J	25 U	9 J	1 U
Toluene	1 10	1 U	1 0	1 0	1 U	25 U	1 U	1 U
Xvienes (total)	<del>                                     </del>	<del></del>	<del>                                     </del>	<del>                                     </del>	<del>                                     </del>	23 0	- ' -	10
Trans-1,2-Dichloroethene	2 U	2 U	2 U	2 U	2 U	50 U	2 U	2 U
Trans-1,3-Dichloropropene	10	10	10	10	1 0	25 U	1 0	1 U
Trichloroethene	2 J	1 0	10	1 0	1 0	25 U		
Trichlorofluoromethane	1 0	10	10	1 0	1 0	25 U	1 U	
Mnyl Chloride	1 0	10	1 U	1 U	1 U			
Anal Culouds		1 0	1 0	<u> </u>	1 1 0	25 U	1 U	1 U

#### Table 5-2 Groundwater Analytical Results - Monitoring Wells - VOCs Evergreen Manor, Roscoe, Illinois

Sample ID:	MW104D-DUP	MW1045	MW105D	MW105D-01	MW105S	MW108D	MW108D-01	MW108D-DUP
Sample Number:	_	_	_	EABX1	_	_	EABW8	
Sample Date:	6/2/00	6/2/00	6/2/00	6/2/00	6/2/00	6/1/00	6/1/00	8/1/00
Laboratory:	ESAT	ESAT	ESAT	MitKem	ESAT	ESAT	MitKem	ESAT
Units	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
PARAMETER	<u> </u>						· · · · · · · · · · · · · · · · · · ·	
1.1.1-Trichloroethane	1 J	2 J	2 J	3	2 J	2 U	1 U	2 U
1,1,2,2-Tetrachioroethane	1 0	1 U	1 0	1 U	1 U	1 U	1 U	1 U
1,1,2-Trichloro-1,2,2-trifluoromethane	2 UJ	2 UJ	2 UJ		2 UJ	2 UJ		2 UJ
1,1,2-Trichloroethane	1 U	1 U	1 U	1 Ü	1 U	1 U	1 U	1 U
1,1-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1.1-Dichloroethene	1 U	1 U	1 U	1 U	1 U	1 Ü	1 0	1 U
1.2.4-Trichlorobenzene		***		1 U		_	1 U	_
1,2-Dibromo-3-chioropropane				1 R			1 R	
1,2-Dibromoethane	1 U	1 U	1 Ü	1 U	1 U	1 U	1 U	1 U
1.2-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 Ü
1,2-Dichloroethane	1 U	1 U	1 0	1 U	1 U	1 U	1 0	1 U
1,2-Dichloropropane	1 0	1 U	1 Ü	1 Ü	1 U	1 Ü	1 0	1 Ü
1,3-Dichlorobenzene	1 U	1 0	1 0	1 0	1 U	1 Ü	1 0	1 Ü
1,4-Dichlorobenzene	1 U	1 U	1 Ü	1 Ü	1 U	1 Ü	1 U	1 Ü
2-Butanone	11 UJ	11 UJ	11 UJ	5 U	11 UJ	11 U	5 U	11 U
2-Hexanone	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U
4-Methyl-2-pentanone	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U
Acetone	2 UJ	2 UJ	2 UJ	5 U	2 UJ	2 U	5 U	2 U
Benzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromochioromethane		***	_	1 U			1 Ü	
Bromodichioromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromomethane	1 U	ט ד	1 U	1 U	1 U	1 U	1 U	1 U
Carbon Disulfide	1 U	1	1 U	1 U	1 U	1 U	1 U	1 U
Carbon Tetrachloride	2 U	2 U	2 U	1 U	2 U	2 U	1 U	2 U
Chiorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroform	3 U	3 U	3 U	1 U	3 U	3 U	1 U	3 U
Chloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 Ü
Cls-1,2-Dichloroethene	4 U	4 U	1 J	2	1 J	4 U	1 U	4 U
Cls-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Dibromochloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Dichlorodifluoromethane	1 U	1 U	1 U		1 U	1 U	-	1 U
Ethylbenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-&/or p-Xylene	1 U	1 U	1 U		1 U	1 U		1 U
Methyl Acetate	1 UJ	1 🖰	1 UJ		1 UJ	1 UJ		1 UJ
Methyl tert-Butyl Ether	1 UJ	1 00	1 UJ		1 UJ	1 U		1 U
Methylene Chloride	1 U	1 U	1 U	2 U	1 U	1 U	2 U	1 0
o-xylene	1 U	1 U	1 U		1 U	1 U		1 U
Styrene	3 UJ	3 W	3 UJ	1 U	3 UJ	3 UJ	1 U	3 UJ
Tetrachloroethene	1 U	1 Ū	3 J	4	3 J	0.6 J	0.7 J	0.6 J
Toluene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Xylenes (total)				1 U			1 U	
Trans-1,2-Dichioroethene	2 U	2 U	2 U	1 U	2 U	2 U	1 U	2 U
Trans-1,3-Dichioropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichioroethene	1 U	1 U	2 J	3	2 J	1 U	1 U	1 U
Trichlorofluoromethane	1 U	1 U	1 U		1 U	1 U		1 U
Mnyl Chloride	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U

Table 5-2
Groundwater Analytical Results - Monitoring Wells - VOCs
Evergreen Manor, Roscoe, Illinois

Sample ID:	MW1085	MW110D	MW110S	MW112
Sample Number:				
Sample Date:	6/1/00	6/1/00	6/1/00	6/2/00
Laboratory:	ESAT	ESAT	ESAT	ESAT
Units	ug/L	ug/L	ug/L	ug/L
PARAMETER				
1,1,1-Trichloroethane	2 U	2 U	2 U	2 J
1,1,2,2-Tetrachioroethane	1 U	1 U	1 U	1 U
1,1,2-Trichloro-1,2,2-trifluoromethane	2 UJ	2 UJ	2 UJ	2 UJ
1,1,2-Trichloroethane	1 U	1 U	1 U	1 U
1,1-Dichloroethane	1 U	1 U	1 U	1 U
1,1-Dichloroethene	1 U	1 U	1 U	1 U
1,2,4-Trichiorobenzene				
1,2-Dibromo-3-chioropropane				
1,2-Dibromoethane	1 U	1 U	1 0	1 U
1,2-Dichiorobenzene	1 U	1 U	1 U	1 U
1,2-Dichioroethane	1 U	1 U	1 U	1 U
1,2-Dichloropropane	1 U	1 U	1 U	1 U
1,3-Dichlorobenzene	1 U	1 U	1 U	1 U
1,4-Dichiorobenzene	1 U	1 U	1 U	1 U
2-Butanone	11 U	11 U	11 U	11 UJ
2-Hexanone	5 U	5 U	5 U	5 U
4-Methyl-2-pentanone	5 U	5 U	5 U	5 U
Acetone	2 U	2 U	2 U_	2 UJ
Benzene	1 U	1 U	1 U	1 U
Bromochloromethane				
Bromodichloromethane	1 U	1 U	1 U	1 U
Bromoform	1 0	10	10	1 U
Bromomethane	1 U	1 U	1 U	1 U
Carbon Disuifide	1 U	1 U	1 U	1 U
Carbon Tetrachioride	2 U	2 U	2 U	2 U
Chlorobenzene	1 U	1 U	1 U	1 U
Chieroethane	1 U	1 U	1 U	1 U
Chloroform	3 U	3 U	3 Ū	3 U
Chloromethane	1 U	1 U	1 U	1 U
Cle-1,2-Dichloroethene	4 U	4 U	4 U	4 U
Cts-1,3-Dichioropropene	1 U	1 U	1 U	1 U
Dibromochioromethane	1 U	1 U	1 U	1 U
Dichlorodifluoromethane	1 U	1 U	1 U	10
Ethylbenzene		1 U		
m-8/or p-Xylene	+	1 113	1 U	1 U
Methyl Acetate			1 UJ	1 UJ
Methyl tert-Butyl Ether	1 U	1 U	1 U	
Methylene Chloride	1 U	1 U	1 U	1 U
o-xylene	1 U 3 UJ	3 UJ	1 U 3 UJ	1 U
Styrene Tetrachioroethene	1 U	1 U		1 U
Toluene	1 0	1 U	1 Ü	1 U
Xylenes (total)	1	<del> </del>	<del></del>	
Trans-1,2-Dichloroethene	2 U	2 U	2 U	2 U
Trans-1,3-Dichloropropene	1 U	1 U	1 U	1 U
Trichloroethene	1 U	1 U	1 U	1 U
Trichlorofluoromethane	1 U	1 U	1 U	1 U
Vinyl Chloride	1 U	1 0	1 U	1 U

U - Constituent not detected; method detection limit (MDL) of the analysis is reported.

J - Concentration reported is an estimated value.

UJ - The analyte was not detected above the reported sample quantitation limit.

DL - Designetes sample was diluted.

B - Designates the constituent was detected in the method blank.

<sup>---</sup> Indicates compound not analyzed.

Evergreen Manor Remedial Investigation Report Section: 5 Revision: 1 Date: 28 March 2001 Page: 25 of 35

# Table 5-3 Groundwater Analytical Results - Monitoring Wells- Water Quality Parameters Evergreen Manor, Roscoe, Illinois

Sample ID:	MW101D-01	MW102D-01	MW103D-01	MW104D-01	MW106D-01	MW108D-01	MW108D-01DP	MW110D-01	MW112-01
Sample Number:	\$07	<b>S11</b>	S08	S15	S14	<b>S12</b>	D12	\$13	S16
Sample Date:	5/30/00	6/1/00	5/30/00	6/2/00	6/2/00	6/1/00	6/1/00	6/1/00	6/2/00
Laboratory:	CHEMTECH	CHEMTECH	CHEMTECH	CHEMTECH	CHEMTECH	CHEMTECH	CHEMTECH	CHEMTECH	CHEMTECH
Units:	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L
Parameter:									
Chemical Oxygen Demand	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	10
Nitrogen, Ammonia	<0.2	<0.2	<0.2	<0.2	<0.2	<0.2	<0.2	<0.2	<0.2
Nitrogen, Nitrite	7.8	<0.05	<0.05	<0.05	<0.05	<0.05	<0.05	<0.05	<0.05
Nitrogen, Nitrate	7.7	5.8	7.5	5.4	4.5	8.3	8.5	6.6	3.3
Ortho Phosphate	0.025	<0.01	0.69	0.016	<0.01	<0.01	0.011	0.014	0.031
Total Sulfide	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Sulfate	27	15	22	25	25	25	22	24	22
				Field Measuremen	ts				
Dissolved Oxygen	6.02	6.95	6.09		2.77	5.57	5.57	5.88	7.61
Ferrous fron	0.0	0.0	0.0		0.0	0.0	0.0	0.0	0.0
Units:	mV	mV	mV	m∨	mV	m∨	m∨	m∨	m∨
Oxidation/Reduction Potential	137	84	28	213	149	174	174	136	167

Table 5-4
Groundwater Analytical Results - Residential Wells - VOCs
Evergreen Manor, Roscoe, Illinois

Sample ID:	RW01-01	RW02-01	RW-03	RW-04	RW04-01	RW-05	RW-05DUP	RW-06
Sample Number:	_	_	_		EABQ4	_		_
Sample Date:	5/31/00	5/31/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00
	12023	12009	11990 Wagon	11990 Blue	11990 Blue			
Address:	Tresemer	Tresemer	Ln.	Spruce	Spruce	4514 Straw Ln.	4514 Straw Ln.	4532 Straw Ln.
Laboratory:	ESAT	ESAT	ESAT	ESAT	MITKEM	ESAT	ESAT	ESAT
Units	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
PARAMETER								
1,1,1-Trichloroethane	2 U	2 U	2 J	2 J	2	2 U	2 U	2 U
1,1,2,2-Tetrachloroethane	1 Ü	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1,2-Trichloro-1,2,2-trifluoromethane	2 UJ	2 UJ	2 U	2 U		2 U	2 U	2 U
1,1,2-Trichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1-Dichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2,4-Trichlorobenzene					1 U			
1,2-Dibromo-3-chloropropane					1 R			
1,2-Dibromoethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichiorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 0
1,2-Dichloroethane	1 U	1 U	1 U	1 U	1 UJ	1 U	1 0	1 U
1,2-Dichloropropane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,3-Dichlorobenzene	1 U					1 U	1 U	1 U
1,A-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U 5 U	1 U	1 U	1 U
2-Butanone 2-Hexanone	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U
	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U
4-Methyl-2-pentanone Acetone	2 U	2 U	0.8	2 U	5 U	2 U	0.6	2 U
Benzese	1 U	1 0	1 U	1 U	1 U	1 U	1 U	1 0
Bromochloromethane					1 U	<u> </u>		<del></del>
Bromodichioromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	1 0	1 U	1 U	1 U	1 U	1 0	1 U	1 0
Bromomethane	1 Ü	10	1 U	1 U	1 U	1 U	1 U	1 0
Carbon Disutfide	1 U	1 U	1 Ū	1 U	1 U	1 U	1 U	1 0
Carbon Tetrachioride	2 U	2 U	2 U	2 U		2 U	2 U	2 Ü
Chlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroform	3 U	3 U	3 U	3 U	1 U	3 U	3 U	3 U
Chloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Cis-1,2-Dichloroethene	4 U	4 U	4 U	1 J	2	4 U	4 U	4 U
Cis-1,3-Dichloropropene	1 U	1 U	1 U	_1 U	1 U	1 U	1 U	1 U
Dibromochloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Dichlorodifluoromethane	1 U	1 U	1 U	1 U		1 U	1 U	1 U
Ethylbenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-&/or p-Xylene	1 U	1 U	1 U	1 U		1 U	1 U	1 U
Methyl Acetate	1 UJ	1 UJ	1 UJ	1 UJ		1 UJ	1 UJ	1 UJ
Methyl tert-Butyl Ether	1 U	1 U	1 U	1 U		1 U	1 U	1 U
Methylene Chloride	1 U	1 U	1 U	1 U	2 U	1 U	1 U	1 U
o-xylene	1 U	1 U	1 0	1 U		1 U	1 U	1 U
Styrene	3 UJ	3 UJ	3 U	3	1 U	3 U	3 U	3 U
Tetrachloroethene	1 U	1 U	1 U	2 J	2	1 U	1 0	1 U
Toluene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Xylenes (total)	2 11	2.0	2.11		1 U	2.11		
Trans-1,2-Dichloroethene	2 U	2 U	2 U	2 U	1 U	2 U	2 U	2 U
Trans-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	1 U	1 0	1 U	5 j	6	1 U	1 U	1 0
Trichlorofluoromethane	1 U	1 U	1 U	1 0	4 11	1 U	1 U	1 U
Vinyl Chloride	1 U	1 U	1 U	1 U	1 U	1 U	1_U	1 U

Page 27 of 35

Table 5-4
Groundwater Analytical Results - Residential Wells - VOCs
Evergreen Manor, Roscoe, Illinois

Sample ID:	RW-07	RW-08	RW-09	RW-10	RW-11	RW11-01	RW11-01DP	RW-11DUP
Sample Number:		_	480		_	EABQ5	EABQ6	-
Sample Date:	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00
	12031 Wagon	11943 Wagon	4134 Valerie	4158 Valerie	4234 Valerie	4234 Valerie	4234 Valerie	4234 Valerie
Address:	Ln. Ct.	Ln.	Dr.	Dr.	Dr.	Dr.	Dr.	Dr.
Laboratory:	ESAT	ESAT	ESAT	ESAT	ESAT	MITKEM	MITKEM	ESAT
Units	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
PARAMETER	<u> </u>							
1.1.1-Trichloroethane	5 J	0.9 J	2 U	2 U	1 J	1	1	1 J
1.1.2.2-Tetrachioroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1,2-Trichloro-1,2,2-triffuoromethane	2 U	2 U	2 U	2 U	2 U	_	-	2 U
1.1.2-Trichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1.1-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1-Dichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2,4-Trichiorobenzene		_	_		_	1 U	1 U	
1,2-Dibromo-3-chloropropane	<b>—</b>				<u> </u>	1 R	1 R	
1,2-Dibromoethane	1 υ	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 UJ	1 U	1 U
1,2-Dichloropropane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,3-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,4-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Butanone	11 U	11 U	11 U	11 U	11 U	5 U	5 U	11 U
2-Hexanone	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U
4-Methyl-2-pentanone	5 U	5 Ü	5 U	5 U	5 U	5 U	5 U	5 U
Acetone	2 U	2 U	2 Ú	2 U	2 U	5 U	5 U	2 U
Benzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 Ü
Bromochioromethane	<del>  _ </del>	_		_		1 U	1 U	
Bromodichloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromomethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Carbon Disulfide	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Carbon Tetrachioride	2 U	2 U	2 U	2 U	2 U		_	2 U
Chlorobenzene	1 U	1 U ·	1 U	1 U	1 U	1 U	1 U	1 U
Chloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroform	3 U	0.9 J	3 U	3 U	3 U	1 U	1 U	3 U
Chloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Cis-1,2-Dichloroethene	4 U	4 U	4 U	4 U	4 U	1 U	1 U	4 U
Cis-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Dibromochloromethane	1 U	1 Ú	1 U	1 U	1 U	1 U	1 U	1 U
Dichlorodifluoromethane	1 U	1 U	1 U	1 U	1 U			1 U
Ethylbenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-8/or p-Xylene	1 U	1 U	1 Ü	1 U	1 U		-	1 U
Methyl Acetate	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ			1 UJ
Methyl tert-Butyl Ether	1 U	1 U	1 U	1 U	1 U			1 U
Methylene Chloride	1 U	1 U	1 U	1 U	1 U	2 U	2 U	1 U
o-xylene	1 U	1 U	1 U	1 U	1 U			1 U
Styrene	3 U	3 U	3 U	3 U	3 U	1 U	1 U	3 U
Tetrachloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Xylenes (total)						1 U	1 U	
Trans-1,2-Dichloroethene	2 U	2 U	2 U	2 U	2 U	1 U	1 U	2 U
Trans-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	0.7 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichlorofluoromethane	1 U	1 U	1 U	1 U	1 U			1 U
Vinyl Chloride	1 U	1 Ü	1 U	1 U	1 U	1 U	1 U	1 U

Table 5-4
Groundwater Analytical Results - Residential Wells - VOCs
Evergreen Manor, Roscoe, Illinois

Sample ID:	RW-12	RW-13	RW-14	RW-15	RW-16	RW-17	RW-18	RW-19
Sample Number:	_			_	_			
Sample Date:	6/5/00	6/6/00	6/6/00	6/6/06	6/6/00	6/6/00	6/6/00	E/E/00
Address:	4176 Valerie Dr.		4628 Straw Ln.	4606 Straw Ln.	4570 Straw Ln.	4246 Hononegah	4232 Hononegah	11974 Blue Spruce Dr.
Laboratory:	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT
Units	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
PARAMETER								
1.1.1-Trichloroethane	2 U	2 U	2 U	2 U	2 U	2 U	2 U	1 J
1.1.2.2-Tetrachloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1.1.2-Trichloro-1.2.2-triffuoromethane	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
1.1.2-Trichloroethane	1 U	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ
1,1-Dichloroethane	1 U	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ
1,1-Dichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2,4-Trichlorobenzene	_			_	_			
1,2-Dibromo-3-chloropropane								
1,2-Dibromoethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,2-Dichloropropane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,3-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,4-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U	_1 U	1 U
2-Butanone	11 U	11 UJ	11 UJ	11 UJ	11 UJ	11 UJ	11 UJ	11 UJ
2-Hexanone	5 υ	5 บม	5 UJ	5 บม	5 บม	5 บม	5 UJ	5 UJ
4-Methyl-2-pentanone	5 U	5 U	5 U	5 U	5 U	5 U	5 Ú	5 U
Acetone	2 U	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ
Benzene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromochloromethane								
Bromodichloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 Ü
Bromomethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Carbon Disuffide	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Carbon Tetrachloride	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 Ú
Chlorobenzene	1 U	1 U	1 U	1 U	1 0	1 U	1 U	1 U
Chloroethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Chloroform	3 U	3 U	3 U	3 U	3 U	3 U	3 U	3 U
Chloromethane	1 U	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ
Cis-1,2-Dichloroethene	4 U	4 U	4 U	4 U	4 U	4 U	4 U	4 U
Cis-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Dibromochloromethane	1 0	1 0	1 U	1 U	1 U	1 U	1 U	1 U
Dichlorodifluoromethane	1 0	1 0	1 0	1 0	1 0	1 U	1 U	1 0
Ethylbenzene	1 0	1 U	1 U	1 0	1 U		1 U	1 U
m-8/or p-Xylene	1 11	1 UJ	1 UJ	1 UJ	1 UJ	1 U 1 UJ	1 U	1 U
Methyl Acetate	1 U	1 U	1 U	1 U	1 U	1 UJ		1 UJ
Methyl tert-Butyl Ether Methylene Chloride	1 0	1 0	1 U	1 0	1 U	1 U	1 U	1 U
o-xylene Chioride	1 0	1 0	1 U	10	1 U	1 U	1 U	1 U
Styrene	3 U	3 U	3 U	3 U	3 U	3 U	3 U	3 U
Tetrachloroethene	1 0	1 U	1 0	10	1 U	1 U	1 U	0.9 J
Toluene	1 0	2 B	2 B	2 B	10	2 B	2 B	1 B
Xylenes (total)								<u>-</u>
Trans-1,2-Dichloroethene	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
Trans-1,3-Dichloropropene	1 0	1 0	1 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	1 0	10	1 0	1 0	1 U	1 0	1 U	4 J
Trichlorofluoromethane	10	10	1 0	1 0	1 U	1 0	1 U	1 U
Vinyl Chloride	1 0	10	10	10	1 0	1 0	1 0	
Tinyi onoite			<u> </u>	<u> </u>		, U	1 0	1 U

#### Table 5-4 Groundwater Analytical Results - Residential Wells - VOCs Evergreen Manor, Roscoe, Illinois

Sample Number:		RW-19DUP	RW-20	RW-21	RW-22
Sample Date:   Series   Seri	Sample ID:	KW-1300F	RW-ZU	KVV-21	KW-22
Moderation   Springer Dr.   Tressmer   Tre					-
Address:   Sprues Dr.   Tresemer   Tresemer   Tresemer   Laboratory:   E&AT   E&AT   E&AT   E&AT   E&AT   Ug/L	Sample Date:				
Laboratory:	Address:				
Units	1 -h	<del></del>			
1,1,1-Trichloroethane		+			
1,1,2-Trichloroethane			ugr	ugre	
1.1,2-Tetrachloroethane		06.1	2 11	2 11	211
1,1,2-Trichlororethane					
1 UJ					
1.1-Dichloroethane	<u> </u>				
1.1-Dichloroethene					
1,2,4-Trichlorobenzene					
1,2-Dibromo-3-chloropropane   1,2-Dibromoethane   1 U		<del>- '-'</del>			
1.2-Dichlorobenzene		<del>                                     </del>	<del></del>		
1.2-Dichlorobenzene		1 (1	1 11	1 11	1 U
1.2-Dichloroethane		<u> </u>			
1		<del> </del>			
1					
1					
2-Butanone					
2-Hexanone		<del></del>			-
Amethyl-2-pentanone					
Acetone					
Benzene					
Bromochloromethane					
Decision   Decision	<del></del>				
1 U		1 (1	1 11	1 U	1 U
Description					
Carbon Disulfide         1 U					
Carbon Tetrachloride         2 U		1 U	1 U	1 U	1 U
Chloroethane         1 U <t< th=""><td>Carbon Tetrachioride</td><td>2 U</td><td>2 U</td><td>2 U</td><td>2 U</td></t<>	Carbon Tetrachioride	2 U	2 U	2 U	2 U
Chloroform	Chlorobenzene	1 U	1 U	1 U	1 U
Chloromethane	Chloroethane	1 U	1 U	1 U	1 U
Cis-1,2-Dichloroethene         4 U         1 U	Chloroform	3 U	3 U	3 U	3 U
1 U	Chloromethane	1 UJ	1 UJ	1 UJ	1 UJ
Distribution   Dist	Cis-1,2-Dichloroethene	4 U	4 U	4 U	4 U
Dichlorodifluoromethane	Cis-1,3-Dichloropropene	1 U	1 U	1 U	1 U
Dichlorodifluoromethane	Dibromochloromethane	1 U	1 U	1 U	1 U
Methyl Acetate	Dichlorodifluoromethane	1 U	1 U	1 U	1 U
Methyl Acetate         1 UJ	Ethylbenzene	1 U	1 U	1 U	1 U
Methyl tert-Butyl Ether         1 U	m-&/or p-Xylene	1 U	1 U	1 U	1 U
Methylene Chloride         1 U	Methyl Acetate	1 UJ	1 UJ	1 UJ	1 UJ
Description	Methyl tert-Butyl Ether	1 U	1 U	1 U	
Styrene	Methylene Chloride	1 U	1 U	1 U	1 U
Tetrachloroethene         0.9 J         1 U         1 U         1 U           Toluene         2 B         1 B         1 B         2 B           Xylenes (total)         —         —         —         —           Trans-1,2-Dichloroethene         2 U         2 U         2 U         2 U         2 U           Trans-1,3-Dichloroethene         1 U	o-xylene				
Toluene	Styrene				
Trans-1,2-Dichloroethene   2 U   2 U   2 U   2 U   2 U   2 U   2 U   Trans-1,3-Dichloropropene   1 U	Tetrachioroethene			<del></del>	
Trans-1,2-Dichloroethene         2 U         2 U         2 U         2 U         2 U           Trans-1,3-Dichloropropene         1 U	Toluene	2 B	1 B	1 B	2 B
Trans-1,3-Dichloropropene         1 U <td>Xylenes (total)</td> <td></td> <td></td> <td>1</td> <td></td>	Xylenes (total)			1	
Trichloroethene         4 J         1 U         1 U         1 U           Trichlorofluoromethane         1 U         1 U         1 U         1 U	Trans-1,2-Dichloroethene				
Trichlorofluoromethane 1 U 1 U 1 U 1 U	Trans-1,3-Dichloropropene			<del> </del>	
	Trichloroethene		1 U	1 U	1 U
Vinyl Chloride 1 U 1 U 1 U 1 U	Trichlorofluoromethane		1 U		
	Vinyl Chloride	1 U	1 U	1 U	1 U

U - Constituent not detected; method detection limit (MDL) of the analysis reported.

J - Concentration reported is an estimated value.

U.J. - The analyte was not detected above the reported sample quantition limit.

B - Designates the constituent was detected in the method blank.

<sup>---</sup> Indicates compound not analyzed.

> Section: 5 Revision: 1

Date: 28 March 2001 Page 30 of 35

Table 5-5

#### Groundwater Analytical Results - Residential Wells - Water Quality Parameters Evergreen Manor, Roscoe, Illinois

Sample ID:	RW01-01	RW01-01DP	RW02-01	RW03-01	RW04-01	RW05-01	RW07-01
Sample Number:	S10	D10	S09	S17	S18	S19	S20
Sample Date:	5/31/00	5/31/00	5/31/00	6/5/00	6/5/00	6/5/00	6/5/00
Laboratory:	CHEMTECH	CHEMTECH	CHEMTECH	CHEMTECH	CHEMTECH	CHEMTECH	CHEMTECH
Units:	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L
Parameter:							
Chemical Oxygen Demand	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0
Nitrogen, Ammonia	<0.2	<0.2	<0.2	<0.2	<0.2	<0.2	<0.2
Nitrogen, Nitrite	<0.05	<0.05	<0.05	<0.05	<0.05	<0.05	<0.05
Nitrogen, Nitrate	4.4	4.2	3.9	6.2	6.3	4.4	5.5
Ortho Phosphate	0.048	0.051	0.048	0.051	0.031	0.038	0.029
Total Sulfide	<1.0	<1.0	<1.0	2.4	1.6	<1.0	<1.0
Sulfate	22	27	24	22	19	26	28
		Field	Measurements				
Dissolved Oxygen	7.03	7.03	7.17	7.70	5.36	5.60	5.39
Ferrous Iron	0.0	0.0	0.0	0.0	0.0	0.0	0.0
Units:	mV	mV	mV	mV	mV	mV	mV
Oxidation/Reduction Potential	88	88	88	222	209	165	176

CHLANO1\WP\RAC\036\29672T5-5.XLS

RFW036-2A-AHVH

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# Table 5-6 Sediment Analytical Results - VOCs Evergreen Manor, Roscoe, Illinois

Sample ID:	SD01-01	SD01-01DUP	SD02-01	SD03-01	SD04-01	SD04-01	\$D04-01DP
<u> </u>	3001-01	0001-01001	GD02-01			EABQ9	EABR1
Sample Number:	5/24/00	5/24/00	5/23/00	5/23/00	5/24/00	5/24/00	5/24/00
Sample Date:	ESAT	BBAT	ESAT	ESAT	ESAT	DATACHEM	DATACHEM
Laboratory: Units	+		ug/kg	ug/kg	ug/kg	ug/kg	ug/kg
	ug/kg	ug/kg	opvy	- op.s	- ug/kg	ugray	- upry
PARAMETER	25 U	12 U	12 U				
1,1,1-Trichioroethane	5 U			5 U	5 U	12 U	12 U
1,1,2,2-Tetrachioroethane		5 U	5 U	5 U	5 U	12 U	12 U
1,1,2-Trichloro-1,2,2-trifluoromethane	5 U	5 U	5 U	5 U	5 U	12 U	12 U
1,1,2-Trichloroethane	10 UJ	12 U	12 U				
1,1-Dichloroethane	5 U	5 U	5 U	5 U	5 U	12 U	12 U
1,1-Dichloroethene							12 0
1,2,4-trichlorobenzene	<del> </del>					12 U	12 U
1,2-Dibromo-3-chloropropane	5 U	5 U	5 U	5 U	5 U	12 UJ	12 UJ
1,2-Dibromoethane	5 U	5 U	5 U	5 U	5 U	12 U	12 U
1,2-Dichlorobenzene	5 U	5 U	5 U	5 U	5 U	12 U	12 U
1,2-Dichloroethane	5 U	5 U	5 U	5 U	5 U	12 U	12 U
1,2-Dichloropropane	5 U	5 U	5 U	5 U	5 U	12 U	12 U
1,3-Dichlorobenzene	5 U	5 U	5 U	5 U	5 U	12 U	12 U
1,4-Dichlorobenzene	<del></del>			25 U	25 U	12 U	12 U
2-Butanone	25 U	25 U 25 U	25 U 25 U	25 U	25 U	12 U	12 U
2-Hexanone	25 U			25 U	25 U	12 U	12 U
4-Methyl-2-pentanone	25 U	25 U	25 U 25 U	25 U	25 U	12 U	12 U
Acetone	5 U	25 U 5 U	5 U	5 U	5 U	12 U	2 J
Benzene	5 U	5 U	5 U	5 U	5 U	12 U	12 U
Bromodichloromethane	5 U	5 U	5 U	5 U	5 U	12 U	12 U
Bromoform	20 UJ	20 UJ	20 UJ		20 UJ	12 U	12 U
Bromomethane	5 U			20 UJ 5 U	5 U	12 U	12 U
Carbon Disulfide		5 U	5 U	5 U	5 U	12 U	12 U
Carbon Tetrachloride	5 U		5 U	5 U		12 U	12 U
Chlorobenzene	5 U	5 U				12 0	12 U
Chloroethane	10 U	12 U	42.11				
Chloroform	8 J 5 U	50 UJ 5 U	50 UJ 5 U	50 UJ 5 U	50 UJ 5 U	12 U	12 U
Chloromethane							<del></del>
Cis-1,2-Dichloroethene	25 U	25 U 5 U	25 U	25 U 5 U	25 U 5 U	12 U 12 U	12 U 12 U
Cis-1,3-Dichloropropene			5 U		3 0	12 U	12 U
Cyclohexane	5 U	5 U	5 U	5 U	<u></u>		
Dibromochloromethane	5 U	5 U	5 U	5 U	5 U	12 U	12 U
Dichiorodifiuoromethane	.5 U	5 U	5 U	5 U	5 U	12 U	12 U
Ethylbenzene	.5 0		3 0	3.0	30	12 U	12 U
isopropylbenzene m-&/or p-Xylene	5 U	5 U		5 U	5 U	12 0	12 0
Methyl Acetate	5 J	15 U	9 J	15 U	15 U	12 U	12 U
Methyl tert-Butyl Ether	25 U	12 U	12 U				
Methylcyclohexane	25 0	25 0			<u> </u>	12 U	12 U
Methylene Chloride	10 U	12 U	12 U				
o-xylene	5 U	5 U	5 U	5 U	5 U		
Styrene	5 U	5 U	5 U	5 U	5 U	12 U	12 U
Tetrachloroethene	5 U	5 U	5 U	5 U	5 U	12 U	12 U
Toluene	5 U	5 U	5 U	5 U	5 U		
Trans-1,2-Dichloroethene	50 U	12 U 12 U	0.7 J				
Trans-1,3-Dichloropropene	5 U	5 U	5 U	50 U	5 U		12 U
Trichloroethene	5 U	5 U	5 U	5 U		12 U	12 U
	5 U	5 U	5 U	5 U	5 U	12 U	12 U
Trichlorofluoromethane					5 U	12 U	12 U
Vinyl Chloride	15 UJ	12 U	12 U				
Xylenes (total)						12 U	12 U

#### Table 5-6 Sediment Analytical Results - VOCs Evergreen Manor, Roscoe, Illinois

Sample ID:	SD05-01	SD06-01
Sample Number: Sample Date:	5/24/00	5/23/00
Laboratory:	ESAT	ESAT
Units	ug/kg	ug/kg
PARAMETER		
1,1,1-Trichloroethane	25 U	25 U
1,1,2,2-Tetrachloroethane	5 U	5 U
1,1,2-Trichloro-1,2,2-triffuoromethane	5 Ü	5 U
1,1,2-Trichloroethane	5 U	5 U
1.1-Dichloroethane	10 UJ	10 UJ
1.1-Dichloroethene	5 Ú	5 U
1,2,4-trichlorobenzene	-	
1,2-Dibromo-3-chloropropane		
1,2-Dibromoethane	5 U	5 U
1.2-Dichlorobenzene	5 U	5 U
1,2-Dichloroethane	5 U	5 U
1,2-Dichloropropane	5 U	5 U
1,3-Dichlorobenzene	5 U	5 U
1,4-Dichlorobenzene	5 U	5 U
2-Butanone	25 U	25 U
2-Hexanone	25 U	25 U
4-Methyl-2-pentanone	25 U	25 U
Acetone	25 U	25 U
Benzene	5 U	5 U
Bromodichloromethane	5 U	5 U
Bromoform	5 U	5 U
Bromomethane	20 UJ	20 UJ
Carbon Disulfide	5 U	5 U
Carbon Tetrachioride	5 U	5 U
Chlorobenzene	5 U	5 U
Chloroethane	10 U	10 U
Chloroform	50 UJ	50 UJ
Chioromethane	5 U	5 U
Cis-1,2-Dichloroethene	25 U	25 U
Cls-1,3-Dichloropropene	5 U	5 U
Cyclohexana		-
Dibromochioromethane	5 U	5 U
Dichlorodifluoromethane	5 U	5 U
Ethylbenzene	5 U	5 U
Isopropylbenzene		
m-&/or p-Xylene	5 U	5 U
Methyl Acetate	9 J	15 U
Methyl tert-Butyl Ether	25 U	25 U
Methylcyclohexane		
Methylene Chloride	10 U	10 U
o-xylene	5 U	5 U
Styrene	5 U	5 U
Tetrachioroethene	5 U	<u>5 U</u>
Toluene	5 U	5 U
Trans-1,2-Dichloroethene	50 U	50 U
Trans-1,3-Dichioropropene	5 U	5 U
Trichioroethene	5 U	5 U
Trichiorofluoromethane	5 U	5 U
Vinyl Chioride	15 UJ	15 UJ
Xylenes (total)	<u> </u>	

U - Constituent not detected; method detection limit (MDL) of the analysis reported.

J - Concentration reported is an estimated value.

UJ - The analyte was not detected above the reported sample quantition limit.

<sup>---</sup> Indicates compound not analyzed.

Revision: 1 Date: 28 March 2001 Page: 33 of 35

### Table 5-7 Sediment Analytical Results- Total Organic Carbon Evergreen Manor, Roscoe, Illinois

Sample ID:	SD01-01	SD02-01	SD03-01	SD04-01	SD05-01	SD06-01	SD06-01DP
Sample Number:	S03	S02	S01	S06	S05	S04	D01
Sample Date:	5/23/00	5/23/00	5/23/00	5/24/00	5/24/00	5/23/00	5/23/00
Laboratory:	CHEMTECH	CHEMTECH	CHEMTECH	CHEMTECH	CHEMTECH	CHEMTECH	CHEMITECH
Units:	mg/kg						
Parameter:							
Total Organic Carbon	30000	17000	11000	4200	24000	4700	5000

### Table 5-8 Surface Water Analytical Results - VOCs Evergreen Manor, Roscoe, Illinois

Sample ID:	SW01-01	SW01-01DUP	SW02-01	SW03-01	SW04-01	SW04-01
Sample Number:	<del>  _</del>	_			_	EABW2
Sample Date:	5/23/00	5/23/00	5/23/00	5/23/00	5/23/00	5/29/00
Laboratory:	ESAT	ESAT	ESAT	ESAT	ESAT	MitKem
Jnits	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
PARAMETER					<u> </u>	
.1.1-Trichloroethane	2 U	2 U	2 U	2 U	2 U	1 U
1,2,2-Tetrachioroethane	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ	1 U
,1,2-Trichloro-1,2,2-trifluoromethane	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ	
1,2-Trichloroethane	1 U	1 U	1 U	1 U	1 U	1 U
,1-Dichloroethane	1 U	1 U	1 U	1 U	1 U	1 U
,1-Dichloroethene	1 UJ	1 UJ	1 W	1 UJ	1 UJ	1 U
2.4-trichiorobenzene						1 U
,2-dibromo-3-chloropropane	<del>-</del>		_	***		1 R
,2-Dibromoethane	1 U	1 U	1 U	1 U	1 U	1 U
,2-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U
,2-Dichioroethane	1 U	1 U	1 U	1 U	1 U	1 UJ
,2-Dichloropropane	1 U	1 U	1 U	1 U	1 U	1 U
,3-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U
A-Dichlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U
-Butanone	11 U	11 U	11 U	11 U	11 U	5 U
-Hexanone	5 U	5 U	5 U	5 U	5 U	5 U
-Methyl-2-pentanone	5 U	5 U	5 U	5 U	5 U	5 U
Acetone	2 UJ	2 UJ	2 W	2 UJ	2 W	5 U
lenzene	1 U	1 U	1 U	1 U	1 U	1 U
romodichloromethane	1 U	1 U	1 U	1 U	1 U	1 U
Iromoform	1 U	1 U	1 Ü	1 U	1 U	1 U
Iromomethane	1 U	1 U	1 U	1 U	1 U	1 U
arbon Disulfide	1 U	1 U	1 U	1 U	1 U	1 U
Carbon Tetrachloride	2 U	2 U	2 U	2 U	2 U	1 U
hlorobenzene	1 U	1 U	1 U	1 U	1 U	1 U
Chloroethane	1 U	1 U	1 U	1 U	1 U	1 U
Chloroform	3 U	3 U	3 U	3 U	3 U	1 U
Chloromethane	1 U	1 U	1 U	1 U	1 U	1 U
Cis-1,2-Dichloroethene	4 U	4 U	4 U	4 U	4 U	1 U
Cis-1,3-Dichloropropene	1 U	1 U	1 U	1 U	1 U	1 U
Dibromochloromethane	1 U	1 U	1 U	1 U	1 U	1 U
Dichlorodifluoromethane	1 U	1 U	1 U	1 U	1 U	
Ethylbenzene	1 U	1 U	1 U	1 U	1 U	1 U
n-&/or p-Xylene	1 U	1 U	1 U	1 U	1 U	
Methyl Acetata	1 UJ	1 UJ	1 W	1 UJ	1 UJ	
Nethyl tert-Butyl Ether	1 U	1 U	1 U	1 U	1 U	
Methylene Chloride	1 U	1 U	1 U	<b>1</b> U	1 U	2 U
-xylene	1 U	1 U	1 U	1 U	1 U	
Styrene	3 U	3 U	3 U	3 U	3 U	1 U
etrachloroethene	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	1 U	1 U	1 U	1 U	1 U	1 U
rans-1,2-Dichloroethene	2 U	2 U	2 U	2 U	2 U	1 U
rans-1,3-Dichioropropene	1 U	1 υ	1 U	1 U	1 U	1 U
Trichloroethene	1 UJ	1 UJ	1 UJ	1 UJ	1 UJ	1 U
richiorofluoromethane	1 U	1 U	1 U	1 U	1 U	
/Inyl Chloride	1 U	1 U	1 U	1 U	1 U	1 U
(ylenes (total)						1 U

3.5

### Table 5-8 Surface Water Analytical Results - VOCs Evergreen Manor, Roscoe, Illinois

Sample ID:	SW04-01DP	SW06-01
Sample Number:	EABW1	_
Sample Date:	5/29/00	5/23/00
Laboratory:	MitKem	ESAT
Units	ug/L	ug/L
PARAMETER		·
1,1,1-Trichloroethane	1 U	2 U
1,1,2,2-Tetrachioroethane	1 U	1 UJ
1,1,2-Trichloro-1,2,2-trifluoromethane		2 UJ
1,1,2-Trichloroethane	1 U	1 U
1.1-Dichloroethane	1 U	1 U
1,1-Dichloroethene	1 U	1 UJ
1,2,4-trichlorobenzene	1 U	
1,2-dibromo-3-chloropropane	1 R	
1,2-Dibromoethane	1 U	1 U
1.2-Dichlorobenzene	1 U	1 U
1.2-Dichloroethane	1 UJ	1 U
1,2-Dichloropropane	1 U	1 U
1,3-Dichlorobenzene	1 U	1 U
1,4-Dichlorobenzene	1 U	1 U
2-Butanone	5 U	11 U
2-Hexanone	5 U	5 U
4-Methyl-2-pentanone	5 U	5 U
Acetone	5 U	2 UJ
Benzene	1 U	1 U
Bromodichloromethane	1 U	1 U
Bromoform	1 U	1 U
Bromomethane	1 U	1 U
Carbon Disulfide	1 U	1 U
Carbon Tetrachloride	1 U	2 U
Chlorobenzene	1 U	1 U
Chloroethane	1 U	1 U
Chloroform	1 U	3 U
Chloromethane	1 U	1 U
Cis-1,2-Dichloroethene	1 U	4 U
Cis-1,3-Dichloropropene	1 U	1 U
Dibromochloromethane	1 U	1 U
Dichlorodifluoromethane		1 U
Ethylbenzene	1 U	1 U
m-&/or p-Xylene		1 U
Methyl Acetate		1 UJ
Methyl tert-Butyl Ether	<del>-</del>	1 U
Methylene Chloride	2 U	1 U
o-xylene	<del></del>	1 U
Styrene	1 U	3 U
Tetrachioroethene	1 U	1 U
Toluene	1 U	1 U
Trans-1,2-Dichloroethene	1 U	2 U
Trans-1,3-Dichloropropene	1 U	1 U
Trichloroethene	1 U	1 UJ
Trichlorofluoromethane	<del></del> _	1 U
Vinyi Chloride	1 U	1 U
Xylenes (total)	1 U	

U - Constituent not detected; method detection limit (MDL) of the analysis is reported.

J - Concentration reported is an estimated value.

R - The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria.

UJ - The analyte was not detected above the reported sample quantitation limit.

<sup>---</sup> Indicates compound not analyzed.

Section: 6 Revisión: 1

Date: 28 March 2001 Page: 1 of 11

**SECTION 6** 

GEOLOGIC AND HYDROGEOLOGIC INVESTIGATION RESULTS

This section presents the site geology, site hydrogeology, and site hydrology based on the RI results,

review of existing documents, and data from previous investigations at the site.

6.1 SITE GEOLOGY

The geology of the site is characterized by surficial deposits of fill and topsoil, overlying

glaciofluvial outwash deposits of sand and gravel. A total of 10 locations were investigated using

CPT. Refusal was reached at a depth of 10 to 15 feet below grade at three of these locations. The

remaining seven locations were advanced to depths greater than approximately 90 feet. The

following sub-section describes the lithology based on the CPT results.

6.1.1 CPT Interpretation

The CPT results included the end-bearing resistance, friction along the side of the probe, friction

ratio, and electrical conductivity. Each of these results is plotted on a CPT log in relation to depth.

The logs from each of the CPT locations are presented in Appendix E. In several instances, shallow

refusal was encountered and subsequent CPT advances at the same location were identified

alphabetically (e.g., CPT07, CPT07A, CPT07B, etc.). Refusal was encountered a number of times,

and at locations CPT07, CPT08, and CPT13, the probe could not be advanced past a depth of 10 to

15 feet. CPT was not conducted at locations CPT09, CPT10, and CPT12 due to their proximity to

nearby locations which were logged for stratigraphy data.

Most of the borings were advanced along roadways and through some thickness of fill or reworked

material. This is seen on the logs as a more clay-rich zone extending to depths from about 1 to 10

CHLAN01\WP\WO\RAC\036\29672S-6.WPD

RFW036-2A-AHVH

Section: 6 Revision: 1

Date: 28 March 2001

Page: 2 of 11

feet below grade. The fill is identified on the logs as having a lower end-bearing resistance (tip

resistance), and a higher friction ratio. Underlying the fill is sand and gravel, with localized zones

containing greater amounts of silt.

Several subtle patterns within the sand and gravel were discernible from the CPT logs. In general,

the uppermost portion of the sand and gravel consisted of a zone with an abundance of dense gravel.

This is based on the tip resistance, which exceeded 300 tons per square foot (tsf) and frequently went

off the scale presented on the logs (>450 tsf). Although there was evidence of concrete fill material

in the vicinity of CPT07, CPT08, and CPT13, it is possible that this gravelly zone was the cause of

the shallow refusal at each of these locations.

Underlying the dense gravelly zone, is a zone that has been interpreted as a clean, well graded sand.

This zone is best seen on the logs from CPT05 and CPT06. At these locations, at depths of 16 to

20 feet below grade, the tip resistance is generally less than 150 tsf and the friction ratio is

approximately 1%. The tip resistance also lacks any sharp peaks that would indicate the presence

of gravels or cobbles. It is also possible that this is a coarsening downwards sequence as evidenced

by a gradual increase in the end-bearing resistance. This zone is either absent or not as well graded

at locations CPT01 through CPT03, which are the southern-most CPT locations at the site.

The remainder of the lithology can be described as sand and gravel, with varying amounts of gravel.

A slight increase in the fines content (probably silt) is apparent at depths below approximately 70

feet below grade. This increase in fines is also associated with an increase in the gravel content.

The above interpretations were correlated with the available boring logs from the IEPA installed

monitoring wells. Transects of three cross sections are shown on Figure 6-1, and the cross sections

showing the interpreted correlation through the site are presented as Figures 6-2 through 6-4. These

cross sections show that the subsurface consists predominantly of sand. Variations in the subsurface

CHLAN01\WP\WO\RAC\036\29672S-6.WPD

RFW036-2A-AHVH

Section: 6 Revision: 1

Date: 28 March 2001

Page: 3 of 11

are due to the amount of gravel present, by the presence of localized silt seams, and by the relatively

higher silt content at certain depths. Low permeability silt/clay layers were not encountered in the

subsurface.

6.2 SITE HYDROGEOLOGY

The hydrogeology at the Evergreen Manor site is that of an extensive unconfined sand and gravel

outwash aquifer. Groundwater elevations were found to be consistently the same in shallow and

deep clustered wells, and varied in elevation between 722 and 735 feet above MSL. Groundwater

flow at the site is from the northeast to the southwest toward the Rock River, as shown on Figure

6-5. The Rock River is presumed to be the groundwater discharge location. The gradient across the

site is fairly uniform and based on the contours shown on Figure 6-5 is approximately 0.0015 ft/ft.

Hydraulic conductivity test results were conducted in the 1980's by the Illinois Department of

Energy and Natural Resources. Pressure tests were conducted at four well clusters, at depths

between 40 and 80 feet below grade, and an average hydraulic conductivity of 3.8 x 10<sup>-2</sup> cm/sec was

found (Wehrmann, 1984).

Using hydraulic conductivity estimates, groundwater gradient, and effective porosity estimates, an

average linear flow velocity for groundwater in the shallow unconfined alluvial aquifer was

estimated in accordance with Darcy's Law as follows:

 $v = \frac{Ki}{n_*}$ 

Where:

v = Linear groundwater seepage velocity (cm/sec)

K = Hydraulic conductivity (cm/sec)

Section: 6 Revision: 1

Date: 28 March 2001

Page: 4 of 11

i = Horizontal hydraulic gradient (ft/ft)

 $n_e$  = Effective porosity

Using a hydraulic conductivity value of 3.8 x 10<sup>-2</sup> cm/sec, a gradient of 0.0015 ft/ft, and an estimated

effective porosity of 30% for sand and gravel mixtures (Fetter, 1994), an average linear groundwater

flow velocity of 1.9 x 10<sup>-4</sup> cm/sec (0.54 ft/day) was estimated.

**6.3 SITE HYDROLOGY** 

The surface water bodies on the site include Dry Creek and the Rock River. Dry Creek is the only

drainageway that traverses a portion of the site and ultimately flows into the Rock River. The Rock

River is the southern boundary of the site. Because of the permeable nature of the sand and gravel

outwash deposits underlying the site, most of the precipitation is expected to infiltrate into the

subsurface and percolate to the groundwater table. However, Dry Creek will also receive surface

runoff during wet periods, when rainwater ponds, or during heavy rainfall. This investigation

occurred during a relatively wet period and Dry Creek was observed to be approaching bank-full

conditions.

The staff gauge reading from 6 June 2000, indicated that the water level in Dry Creek was

approximately 11 ft higher than the water table elevation in the closest wells (MW110S and D).

Based on these readings, Dry Creek is not expected to be in direct hydraulic connection with the

groundwater table at that location. Dry Creek would be classified as a losing stream at the time of

this investigation, indicating that it would contribute water to the subsurface. The amount of water

that is lost from Dry Creek to the subsurface could be calculated by testing or approximating the

permeability of the channel bottom sediments. However, since the channel bottom sediments are

clay and silt rich, the amount of loss from the stream is expected to be minimal.

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RFW036-2A-AHVH

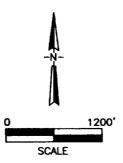
Section: 6 Revision: 1

Date: 28 March 2001

Page: 5 of 11

In the residential areas, primarily south of Hononegah Road, the surface drainage pattern has been somewhat altered by construction of roadways, driveways, and buildings. Although precipitation will percolate through the lawns in the residential area, a portion will be carried by the ditch system to the Rock River.

ELEVATOR RD.



NOTE:
BASEMAP ADAPTED FROM FIGURE 1 OF THE CONTAMINANT SOURCE
EVALUATION — EVERGREEN MANOR SITE, DATED JANUARY 1997, BY
CONESTOGA—ROVERS & ASSOCIATES.

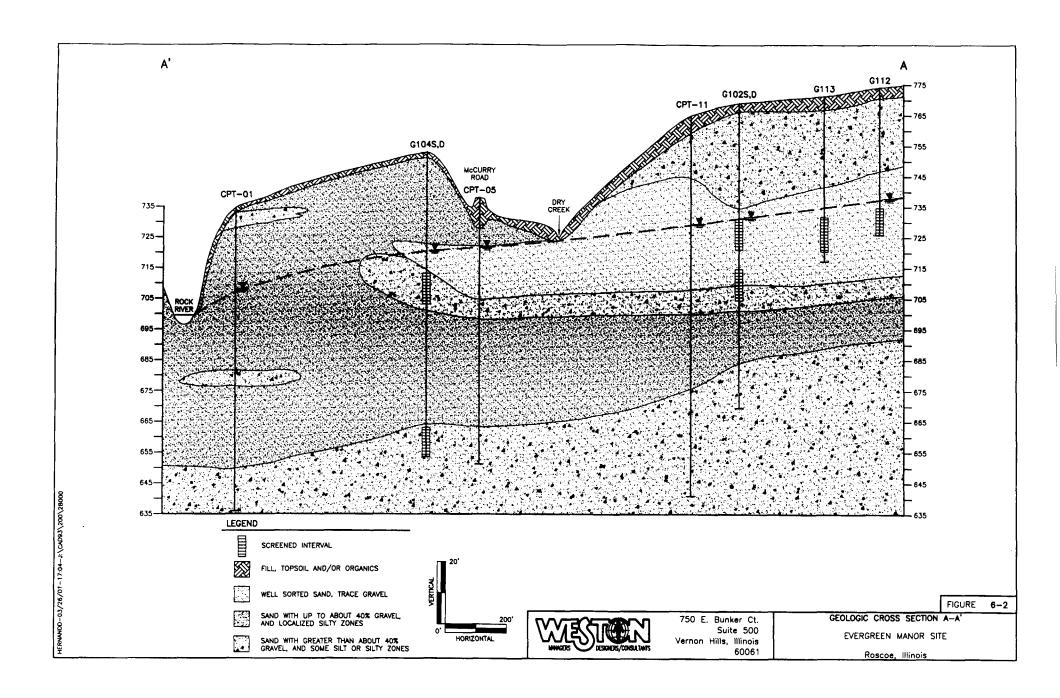
FIGURE 6-1

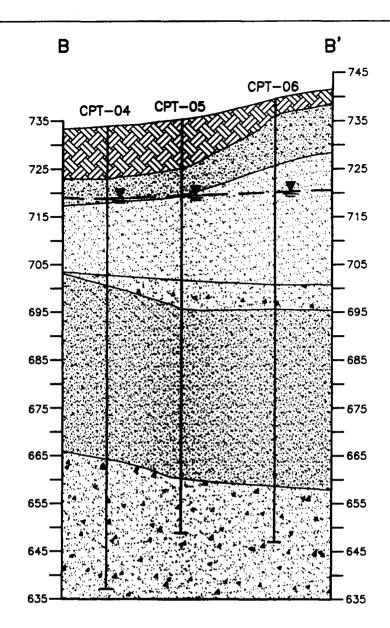


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750 E. Bunker Ct. Suite 500 Vernon Hills, Illinois 60061 CROSS SECTION TRANSECT MAP

EVERGREEN MANOR SITE Roscoe, Illinois





**LEGEND** 



FILL, TOPSOIL AND/OR ORGANICS



HERNANDD-03/26/01-14:36-J:\CAD93\200\28100

WELL SORTED SAND, TRACE GRAVEL



SAND WITH UP TO ABOUT 40% GRAVEL, AND LOCALIZED SILTY ZONES



SAND WITH GREATER THAN ABOUT 40% GRAVEL, AND SOME SILT OR SILTY ZONES

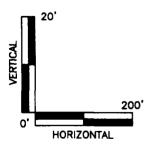


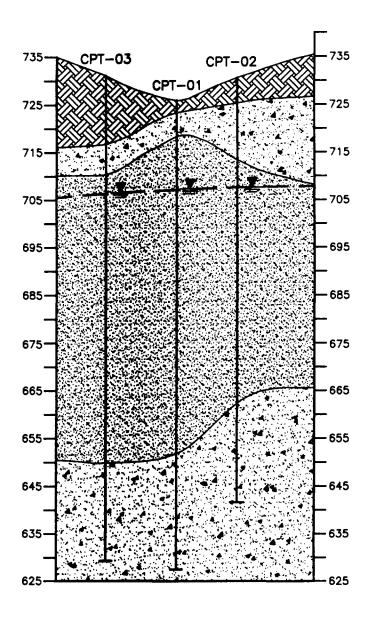
FIGURE 6-3

WANGERS DESIGNERS/CONSULTANTS

750 E. Bunker Ct. Suite 500 Vernon Hills, Illinois 60061 EVERGREEN MANOR SITE

Roscoe, Illinois





**LEGEND** 

FILL, TOPSOIL AND/OR ORGANICS



WELL SORTED SAND, TRACE GRAVEL



SAND WITH UP TO ABOUT 40% GRAVEL, AND LOCALIZED SILTY ZONES



SAND WITH GREATER THAN ABOUT 40% GRAVEL, AND SOME SILT OR SILTY ZONES

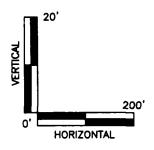


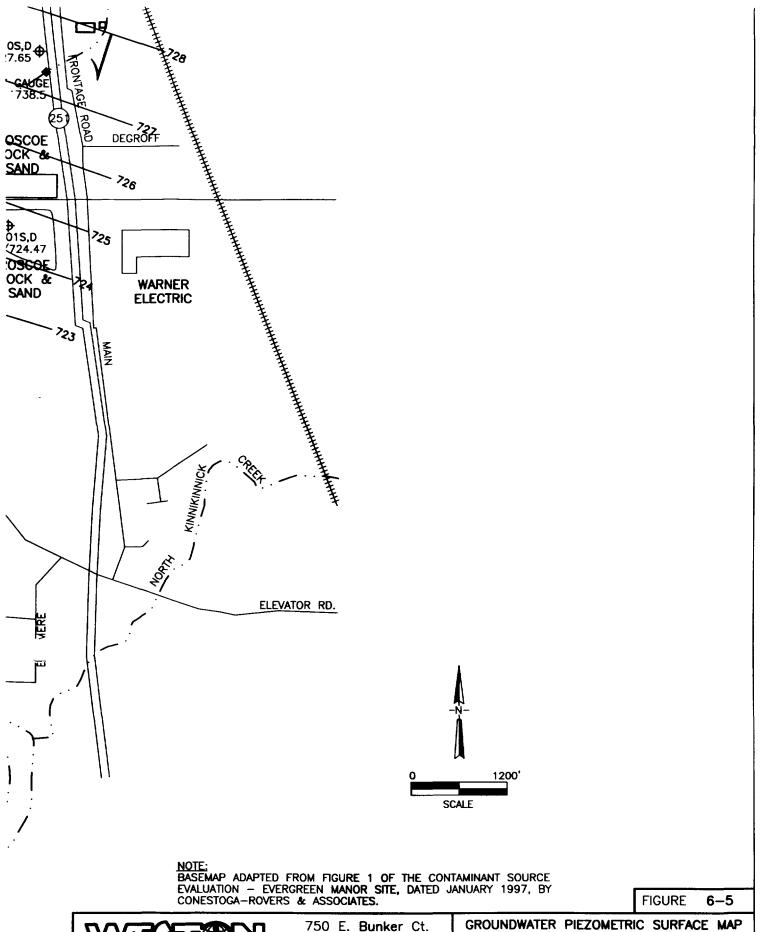
FIGURE 6-4

750 E. Bunker Ct.
Suite 500

Vernon Hills, Illinois 60061 EVERGREEN MANOR SITE Roscoe, Illinois

GEOLOGIC CROSS SECTION C-C'







750 E. Bunker Ct. Suite 500 Vernon Hills, Illinois 60061

GROUNDWATER PIEZOMETRIC SURFACE MAP
6 JUNE 2000
EVERGREEN MANOR SITE
Roscoe, Illinois

> Section: 6 Revision: 1

Date: 28 March 2001

Page: 11 of 11

### TABLE 6-1 **Groundwater Elevations from Monitoring Wells Evergreen Manor** Roscoe, Illinois

	Elevation of Top of Inner	Elevation of Ground	Depth to Bottom (ft)	Total Donth of Wall	Depth to Water (ft)	Groundwater
\A( 1) A D	•		1 ' ' '	•	1 ' ' ' '	0.00
Well I.D.	Casing (ft above MSL)	Surface (ft above MSL	6 June 2000	(ft below grade)	6 June 2000	Elevation
MW-101S	730.29	727.5	42.64	39.9	5.82	724.47
MW-101D	730.34	727.6	82.00	79.3	5.87	724.47
MW-102S	771.09	768.7	47.71	45.3	40.20	730.89
MW-102D	771.49	769.1	67.00	64.6	40.56	730.93
MW-103S	767.21	764.3	45.11	42.2	35.90	731.31
MW-103D	767.27	764.4	58.98	56.1	35.94	731.33
MW-104S	756.13	753.3	62.43	59.6	33.55	722.58
MW-104D	755.59	753.1	102.30	99.8	32.96	722.63
MW-105S	757.95	755.2	67.85	65.1	35.15	722.80
MW-105D	757.79	755.3	101.81	99.3	34.96	722.83
MW-106S	757.20	754.8	67.50	65.1	34.15	723.05
MW-106D	756.86	754.5	102.39	100.0	33.83	723.03
MW-107S	765.79	763.4	47.34	45.0	36.56	729.23
MW-107D	766.39	763.3	67.40	64.3	37.21	729.18
MW-108S	767.01	764.4	46.98	44.4	38.33	728.68
MW-108D	766.96	764.4	67.35	64.8	38.23	728.73
MW-109S	769.91	767.3	52.45	49.8	41.69	728.22
MW-109D	769.50	766.9	72.38	69.8	41.36	728.14
MW-110S	748.19	745.4	32.53	29.7	20.60	727.59
MW-110D	748.31	745.6	52.04	49.3	20.66	727.65
MW-111	770.70		50.65		38.08	732.62
MW-112*	774.2	772.3	50.66	48.8	38.98	735.22

#### NOTES:

\*: Elevations for MW-112 should be considered approximate, since the casing has been bent since the last vertical survey.

---: Information not available

Section: 7 Revision: 1

Date: 28 March 2001 Page: 1 of 20

**SECTION 7** 

NATURE AND EXTENT OF CONTAMINATION

One of the objectives of site characterization activities is to evaluate the nature and extent of

contamination, such that informed decisions can be made regarding the level of risk presented by

the site and the appropriate type of removal action necessary. During this investigation, data were

collected and compared to screening levels to delineate the extent of contamination. The following

subsections discuss the development and applicability of screening levels, potential sources of

contamination, and the nature and extent of contamination in groundwater, surface water, and

sediment.

7.1 <u>DEVELOPMENT OF SCREENING LEVELS</u>

Screening levels were developed to provide a means for determining the extent of contamination at

this site. A screening level is an acceptable level for each contaminant of concern in each exposure

route. The screening levels serve as cleanup levels protective of both human health and the

environment, but are not necessarily used for direct implementation as removal action objectives

(RAOs); however, they can be considered as one factor among many in the development of RAOs.

The primary purpose of the screening levels is to identify constituents of potential concern (COPCs)

for the site and to illustrate the extent of contamination. This section discusses the development of

screening levels for individual contaminants for groundwater, sediment, and surface water.

7.1.1 Groundwater

Groundwater screening levels were developed by evaluating the applicable regulatory standards.

These include the Illinois Environmental Protection Agency (IEPA) groundwater quality standards

and the U.S. EPA groundwater quality standards. The IEPA standards are described in Title 35:

Environmental Protection, Part 742 - Tiered Approach to Corrective Action Objectives (TACO).

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RFW036-2A-AHVH

Section: 7 Revision: 1

Date: 28 March 2001

Page: 2 of 20

Table E, Tier I Groundwater Remediation Objectives for the Groundwater Component of the

Groundwater Ingestion Route, of Appendix B of TACO lists all of the applicable groundwater

quality standards. The U.S. EPA regulatory standards are the Maximum Concentration Levels

(MCLs), which are incorporated into the Safe Drinking Water Act (SDWA) and are found in 40 CFR

141.61 - Maximum Contaminant Levels (MCLs) for Organic Contaminants (Integrated). The most

stringent of the two regulatory standards was chosen as the screening level for each detected

constituent. If a MCL did not exist, the TACO value was used as the screening level. The screening

levels are presented in Table 7-1.

7.1.2 Sediment

The data presented in Section 5.5 indicated that several constituents were detected in sediment. The

screening levels are based on the IEPA standards found in Table A, Tier 1 Soil Remediation

Objectives for Residential Properties, of Appendix B of TACO, and the U.S. EPA Region IX risk-

based concentrations (RBCs) for residential ingestion exposure route (U.S. EPA, 1996). Since

sediment standards do not exist in TACO, the more conservative TACO standards for soil were used.

The most stringent of the two standards was used as the screening level. The screening levels are

presented in Table 7-1 for those compounds which resulted in positive detections during sampling

and analysis.

7.1.3 Surface Water

At a minimum, surface water data would be derived from Ambient Water Quality Criteria.

However, all of the data were reported below method detection limits. Therefore, surface water

screening levels were not developed.

CHLAN01\WP\WO\RAC\036\29672S-7.WPD

RFW036-2A-AHVH

Section: 7 Revision: 1

Date: 28 March 2001

Page: 3 of 20

7.2 SOURCES OF CONTAMINATION

The investigative work performed during this RI, as well as previous investigations, have attempted

to identify potential sources of VOC contamination. The previous investigations were not able to

identify a unique source of contamination; they were only able to generalize by indicating the source

was located in or near the industrial park in the vicinity of Rockton Road and IL Route 251. The

analytical results of this RI indicated that most of the contaminants detected were at low

concentrations, close to the detection limits, and were likewise not able to pinpoint a specific

contaminant source. The fracture trace analysis results indicated that the industrial park near

Rockton Road and IL Route 251 could serve as a source area.

7.3 EXTENT OF CONTAMINATION IN GROUNDWATER

Groundwater samples were collected from 10 CPT locations, 15 monitoring wells, and 22 residential

wells and analyzed for VOCs. A total of 14 distinct VOCs were found in groundwater samples

above the method detection limits. Tables 7-2 through 7-4 present a summary of the detected

constituents, and their respective screening levels, found in groundwater from CPT, monitoring well,

and residential well sampling, respectively. The highlighted values on these tables are the results

which exceeded screening levels.

A total of three compounds were found to exceed screening levels. Trichloroethene was found in

residential well RW04 at a concentration of 6 µg/L. Chloroform was found in residential well RW08

at a concentrations of 0.9 µg/L. Tetrachloroethene was found in monitoring well MW103S at a

concentration of 9 µg/L. No exceedances were found in the CPT samples. Thus, chloroform,

tetrachloroethene, and trichloroethene are considered COPCs at the site.

Figure 7-1 shows the extent of detected VOCs in the vicinity of Evergreen Manor. This figure also

identifies the three locations where screening level exceedances occurred. In large part, the extent

CHLAN01\WP\WO\RAC\036\29672S-7.WPD

RFW036-2A-AHVH

Section: 7 Revision: 1

Date: 28 March 2001

Page: 4 of 20

of the detected VOCs is based on the presence of acetone and toluene, which are present at low

levels but are not considered COPCs.

The extent of contamination in groundwater is considered to occur only locally at the locations

where screening level exceedances occurred. This includes the area near MW103, RW04, and

RW08. These areas are within the area defined previously by other investigations.

7.4 EXTENT OF CONTAMINATION IN SEDIMENT

Table 7-5 shows the constituents detected in sediment and the screening level for each of the

detected constituents. The data in this table indicated that all of these constituents were detected

at concentrations considerably less than their respective screening levels. Therefore, based on the

absence of screening level exceedances, none of the constituents detected in sediment are considered

COPCs, and the sediment at the site is not considered contaminated.

7.5 EXTENT OF CONTAMINATION IN SURFACE WATER

Based on the fact that all of the surface water data were reported as not detected above the method

detection limits, none of the analyzed constituents are considered COPCs, and the surface water is

not considered contaminated.

CHLAN01\WP\WO\RAC\036\29672S-7.WPD

RFW036-2A-AHVH

Section: 7

Revision: 1 Date: 28 March 2001

te: 28 March 2001 Page: 8 of 20

## Table 7-2 Summary of Detected Consituents in Groundwater CPT - VOCs

Sample ID:		CPT-02-02	CPT02-03	CPT-02-03	CPT-02-04	CPT-02-05	CPT-02-06	CPT-02-07	CPT-03-01	CPT-03-02	CPT-03-02DP
Sample Number:			EABX4								
Sample Date:		6/2/00	6/2/00	6/2/00	6/2/00	6/2/00	6/2/00	6/2/00	5/26/00	5/26/00	5/26/00
Laboratory:		ESAT	MitKem	ESAT							
Sample Screen Depth (1	it below ground):	34	42	42	51	68	78	84	29	42	42
Units:		ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Parameter (Screen	ing Level)										
Acetone	(700)	2 UJ	470	55 J	2 UJ	2 UJ	2 UJ	2 UJ	20 B	32 B	19 B
Methylene Chloride	(5)	1 U	10 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1-Dichloroethane	(700)	1 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Butanone	()	11 UJ	25 U	11 UJ	11 U	11 U	11 U				
Cis-1,2-Dichloroethene	(70)	4 U	5 U	4 U	4 U	4 U	4 U	4 U	4 U	4 U	4 U
1,1,1-Trichloroethane	(200)	2 U	5 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
Benzene	(5)	1 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	(5)	1 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	(1000)	0.6 J	5 U	0.7 J	0.5 J	1 J	1 J	1 U	0.8 J	1 J	0.8 J
Tetrachloroethene	(5)	1 U	5 U	1 Ü	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Ethylbenzene	(700)	1 Ü	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-&/or p-Xylene	(10,000)	0.5 J		1 U	1 U	1 U	1 U	0.6 J	1 U	1 U	1 U
o-xylene	(10,000)	1 U		1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Xylenes (total)	(10,000)		5 U					_			

Section: 7 Revision: 1

Date: 28 March 2001 Page: 9 of 20

### Table 7-2 Summary of Detected Consituents in Groundwater CPT - VOCs

#### Evergreen Manor, Roscoe, Illinois

Sample ID:		CPT-03-05	CPT03-06	CPT-03-06	CPT-03-07	CPT-04-01	CPT-04-02	CPT-04-02	CPT04-02DP	CPT-04-02DUP
Sample Number:			EABX5	-	_			EABW5	EABW6	
Sample Date:		5/26/00	6/2/00	6/2/00	6/2/00	5/29/00	5/29/00	5/29/00	5/29/00	5/29/00
Laboratory:		ESAT	MitKem	ESAT	ESAT	ESAT	ESAT	MitKem	MitKem	ESAT
Sample Screen Depth	(ft below ground):	84	92	92	102	32	46	46	46	46
Units:		ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Parameter (Scree	ning Level)									
Acetone	(700)	9 J	5 U	2 UJ	2 UJ	2 U	10 B	5 U	5 U	2 U
Methylene Chloride	(5)	0.5 J	2 U	1 U	1 U	1 U	1 U	2 U	2 U	1 U
1,1-Dichloroethane	(700)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Butanone	()	11 U	5 U	11 UJ	11 UJ	11 U	11 U	5 U	5 U	11 U
Cis-1,2-Dichloroethene	(70)	4 U	1 U	4 U	4 U	4 U	4 U	1 U	1 U	4 U
1,1,1-Trichloroethane	(200)	2 U	0.9 J	0.9 J	0.8 J	2 U	2 U	1 U	1 U	2 U
Benzene	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	(1000)	1 J	2	2 J	2 J	1 J	0.8 J	0.6 J	0.8 J	0.9 J
Tetrachloroethene	(5)	1 U	1 U	1 U	1_U	1 U	1 U	1 U	1 U	1 U
Ethylbenzene	(700)	1 U	1 U	1 U	1	1 U	1 U	1 U	1 U	1 U
m-&/or p-Xylene	(10,000)	1 U		1 U	1 U	1 U	1 U			1 U
o-xylene	(10,000)	1 U		1 U	1 U	1 U	1 U			1 U
Xylenes (total)	(10,000)		1 U	_				1 U	1 U	

CHLANO1\WP\RAC\036\29672T7-2.XLS

RFW036-2A-AHVH

Section: 7 Revision: 1

Date: 28 March 2001 Page: 10 of 20

### Table 7-2 Summary of Detected Consituents in Groundwater CPT - VOCs

Sample ID:		CPT-04-05	CPT-04-06	CPT-04-07	CPT-05-01	CPT-05-02	CPT-05-03	CPT-05-04	CPT-05-05DUP	CPT-05-06	CPT-05-07
Sample Number:					-						
Sample Date:		5/29/00	5/29/00	5/29/00	5/29/00	5/29/00	5/29/00	5/29/00	5/30/00	5/30/00	5/30/00
Laboratory:		ESAT	ESAT	ESAT							
Sample Screen Depth (f	t below ground):	78	84	93	35	43	51	57	69.5	78	87
Units:		ug/L	ug/L	ug∕L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Parameter (Screen	ing Level)										
Acetone	(700)	2 U	8 B	9 B	2 U	2 U	2 U	2 U	2 U	100 J	2 U
Methylene Chloride	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1-Dichloroethane	(700)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Butanone	()	11 U	16 J	11 U							
Cis-1,2-Dichloroethene	(70)	4 U	4 U	4 U	4 U	4 U	4 U	4 U	4 U	4 U	4 U
1,1,1-Trichloroethane	(200)	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U	2 U
Benzene	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	(1000)	1 J	0.8 J	1 J	1 J	1 J	1 J_	1 J	0.6 J	0.8 J	1 J
Tetrachloroethene	(5)	1 U	1 U_	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Ethylbenzene	(700)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-&/or p-Xylene	(10,000)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
o-xylene	(10,000)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Xylenes (total)	(10,000)				-						

Section: 7

Revision: 1 Date: 28 March 2001

ate: 28 March 2001 Page: 11 of 20

### Table 7-2 Summary of Detected Consituents in Groundwater CPT - VOCs

Sample ID:		CPT-06-02	CPT-06-03	CPT-06-04	CPT-06-05	CPT06-06	CPT-06-06	CPT-09-01	CPT-09-04	CPT-09-05	CPT-09-06
Sample Number:						EABW9					
Sample Date:		6/1/00	6/1/00	6/1/00	6/1/00	6/1/00	6/1/00	5/30/00	5/30/00	5/30/00	5/30/00
Laboratory:		ESAT	ESAT	ESAT	ESAT	MitKem	ESAT	ESAT	ESAT	ESAT	ESAT
Sample Screen Depth (	ft below ground):	42	53	62	74	85	85	35	68	75	85
Units:		ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Parameter (Scree	ning Level)										
Acetone	(700)	53 J	11 J	7 J	11 J	7 U	2 U	2 U	2 U	2 U	11 J
Methylene Chloride	(5)	1 U	1 U	1 U	1 U	2 U	1 U	1 U	1 U	1 U	1 U
1,1-Dichloroethane	(700)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Butanone	()	11 U	11 U	11 U	11 U	5 U	11 U				
Cis-1,2-Dichloroethene	(70)	4 U	4 U	4 U	4 U	1 U	4 U	4 U	4 U	4 U	4 U
1,1,1-Trichloroethane	(200)	2 U	2 U	2 U	2 U	1 U	2 U	2 U	2 U	2 U	2 U
Benzene	(5)	1 U	1 U	1 U	1 U	1 U	1 U	0.5 J	1 U	1 U	1 U
Trichloroethene	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	(1000)	0.8 J	0.8 J	1 J	1 J	2	2 J	1 J	0.8 J	0.5 J	0.8 J
Tetrachioroethene	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Ethylbenzene	(700)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-&/or p-Xylene	(10,000)	1 U	1 U	1 U	1 U		1 U	1 U	1 U	1 U	1 U
o-xylene	(10,000)	1 U	1 U	1 U	1 U		1 U	1 U	1 U	1 U	1 U
Xylenes (total)	(10,000)					1 U					

Section: 7 Revision: 1

Date: 28 March 2001

: 28 March 2001 Page: 12 of 20

## Table 7-2 Summary of Detected Consituents in Groundwater CPT - VOCs

Sample ID:		CPT-10-01	CPT-10-03	CPT-10-04	CPT-10-4DUP	CPT-10-05	CPT-10-06	CPT-10-07	CPT11-01	CPT-11-01
Sample Number:									EABX7	
Sample Date:		5/31/00	5/31/00	6/1/00	6/1/00	6/1/00	6/1/00	6/1/00	6/3/00	6/3/00
Laboratory:		ESAT	ESAT	ESAT	ESAT	ESAT	ESAT	ESAT	MitKem	ESAT
Sample Screen Depth (	ft below ground):	25	42	55	55	65	73	90	45	45
Units:		ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L.
Parameter (Screer	ning Level)									
Acetone	(700)	2 U	40 B	2 U	2 U	2 U	24 J	2 U	5 U	2 UJ
Methylene Chloride	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	2 U	1 U
1,1-Dichloroethane	(700)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Butanone	()	11 U	11 U	11 U	11 U	11 U	11 U	11 U	5 U	11 UJ
Cis-1,2-Dichloroethene	(70)	4 U	4 U	4 U	4 U	4 U	4 U	4 U	1 U	4 U
1,1,1-Trichloroethane	(200)	2 U	2 U	1 J	1 J	0.8 J	2 U	2 U	1 U	2 U
Benzene	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	(1000)	0.9 J	1 J	0.5 J	1 U	0.9 J	1 J	1 J	1 U	1 U
Tetrachioroethene	(5)	1 U	1 U	0.6 J	0.6 J	1 U	1 U	1 U	1 U	1 U
Ethylbenzene	(700)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-&/or p-Xylene	(10,000)	1 U	1 U	1 U	1 U	1 U	1 U	1 U		1 U
o-xylene	(10,000)	1 U	1 U	1 U	1 U	1 U	1 U _	1 U		1 U
Xylenes (total)	(10,000)								0.6 J	

Revision: 1

Date: 28 March 2001 Page: 13 of 20

### Table 7-2 **Summary of Detected Consituents in Groundwater CPT - VOCs** Evergreen Manor, Roscoe, Illinois

Sample ID:		CPT-11-04	CPT11-05	CPT-11-05	CPT11-05DP	CPT-11-06	CPT-11-07	CPT11-08	CPT-11-08	CPT-12-01	CPT-12-01DUP
Sample Number:			EABX9		EABQ1			EABQ2			
Sample Date:		6/3/00	6/3/00	6/3/00	6/3/00	6/3/00	6/3/00	6/3/00	6/3/00	6/5/00	6/5/00
Laboratory:		ESAT	MitKem	ESAT	MitKem	ESAT	ESAT	MitKem	ESAT	ESAT	ESAT
Sample Screen Depth (fi	below ground):	81	93	93	93	102	114	125	125	45	
Units:	<u> </u>	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Parameter (Screen	ing Level)										
Acetone	(700)	2 UJ	5 U	2 UJ	5 U	9 J	13 J	5 U	2 UJ	1 B	1 B
Methylene Chloride	(5)	1 U	2 U	1 U	2 U	1 U	1 U	2 U	1 U	1 U	1 U
1,1-Dichloroethane	(700)	1 U	2	1 U	2	1 U	1 U	1 U	1 UJ	1 U	1 U
2-Butanone	()	11 UJ	5 U	11 UJ	5 U	11 UJ	11 UJ	5 U	11 UJ	11 U	11 U
Cis-1,2-Dichloroethene	(70)	4 U	1 U	4 U	1 U	4 U	4 U	1 U	4 U	4 U	4 U
1,1,1-Trichloroethane	(200)	0.8 J	2	2 J	3	3 J	2 U	0.7 J	2 U	2 U	2 U
Benzene	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	(5)	1 U	1 U	1 U	1 U	1 Ü	1 U	1 U	1 U	1 U	1 U
Toluene	(1000)	0.9 J	2	1 J	0.8 J	1 J	2 J	2	3 J	0.5 J	0.5 J
Tetrachloroethene	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Ethylbenzene	(700)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-&/or p-Xylene	(10,000)	1 U		1 U		1 U	1 U		1 U	1 U	1 U
o-xylene	(10,000)	1 U		1 U		1 U	1 U		1 U	1 U	1 U
Xylenes (total)	(10,000)		1 U		1 U			1 U			

Section: 7 Revision: 1

Date: 28 March 2001

Page: 14 of 20

### Table 7-2 Summary of Detected Consituents in Groundwater CPT - VOCs

#### Evergreen Manor, Roscoe, Illinois

Sample ID:		CPT-12-03	CPT-12-04	CPT-12-04DUP	CPT-12-05	CPT12-06	CPT-12-06	CPT-12-07
Sample Number:		EABQ3		_		EABQ7		
Sample Date:		6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00
Laboratory:		MitKem	ESAT	ESAT	ESAT	MitKem	ESAT	ESAT
Sample Screen Depth (ft	below ground):	70	81	81	93	102	102	118
Units:		ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Parameter (Screeni	ng Level)				<del></del>			
Acetone	(700)	5 U	2 B	7 B	2	5 U	2	2
Methylene Chloride	(5)	2 U	1 U	1 U	1 U	2 U	1 U	1 U
1,1-Dichloroethane	(700)	1 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Butanone	()	5 U	11 U	11 U	11 U	5 U	11 U	11 U
Cis-1,2-Dichloroethene	(70)	1 U	4 U	4 U	4 U	1 U	4 U	4 U
1,1,1-Trichloroethane	(200)	1 U	2 U	2 U	2 U	1 U	2 U	2 U
Benzene	(5)	1 U	1 U	0.5 J	1 U	1 U	1 U	1 U
Trichloroethene	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	(1000)	0.8 J	0.9 J	1 U	2 J	1	2 J	2 J
Tetrachloroethene	(5)	1 Ü	1 U	1 U	1 Ú	1 U	1 U	1 U
Ethylbenzene	(700)	1 U	1 U	1 U	1 U	1 U	1 U	1 U
m-&/or p-Xylene	(10,000)		1 U	1 U	1 U		1 U	1 U
o-xylene	(10,000)		1 U	1 U	1 U		1 U	1 U
Xylenes (total)	(10,000)	1 U				1 U		

Highlighting indicates compound exceeded the Screening Level.

- U Constituent not detected; method detection limit (MDL) of the analysis reported.
- J Concentration reported is an estimated value.
- UJ The analyte was not detected above the reported sample quantition limit.
- B Designates the constituent was detected in the method blank.
- -- Indicates compounds not analyzed.

Section: 7 Revision: 1

Date: 28 March 2001

: 28 March 2001 Page: 15 of 20

# Table 7-3 Summary of Detected Constituents in Groundwater Monitoring Wells - VOCs Evergreen Manor, Roscoe, Illinois

Sample ID:	MW101D	MW101S	MW102D	MW102S	MW103D	MW103D-DL	MW103S	MW104D	MW104D-DUP
Sample Number:							•••	1	
Sample Date:	5/30/00	5/30/00 6/1/00		6/1/00	5/31/00	5/31/00	5/31/00	6/2/00	6/2/00
Laboratory:	ESAT	ESAT ESAT		ESAT	ESAT	ESAT	ESAT	ESAT	ESAT
Units	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
PARAMETER (Screening Level)									
1,1,1-Trichloroethane (200)	2 U	2 U	2 U	2 U	3 J	50 U	2 U	1 J	1 J
1,1,2-Trichloro-1,2,2-trifluoromethane ()	2 UJ	2 UJ	<b>2</b> U	2 UJ	300 J	180 J	2 J	2 UJ	2 UJ
Cis-1,2-Dichloroethene (70)	4 U	4 U	4 U	4 U	4 U	100 U	4 U	4 U	4 U
Tetrachloroethene (5)	1 U	1 U	1 U	1 U	0.5 J	25 U		1 U	1 U
Trichloroethene (5)	2 J	1 U	1 U	1 U	1 U	25 U	1 U	1 U	1 U

Section: 7 Revision: 1

Date: 28 March 2001

Page: 16 of 20

## Table 7-3 Summary of Detected Constituents in Groundwater Monitoring Wells - VOCs Evergreen Manor, Roscoe, Illinois

Sample ID:		MW104S	MW105D	MW105D-01	MW105S	MW108D	MW108D-01	MW108D-DUP	MW108S	MW110D	MW110S	MW112
Sample Number:			-	EABX1			EABW8					
Sample Date:		6/2/00	6/2/00	6/2/00	6/2/00	6/1/00	6/1/00	6/1/00	6/1/00	6/1/00	6/1/00	6/2/00
Laboratory:		ESAT	ESAT	MitKem	ESAT	ESAT	MitKem	ESAT	ESAT	ESAT	ESAT	ESAT
Units		ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
PARAMETER (Screenin	g Level)											
1,1,1-Trichloroethane	(200)	2 J	2 J	3.00	2 J	2 U	1 U	2 U	2 U	2 U	2 U	2 J
1,1,2-Trichloro-1,2,2-trifluorometh	ane ()	2 UJ	2 UJ		2 UJ	2 UJ		2 UJ	2 UJ	2 UJ	2 UJ	2 UJ
Cis-1,2-Dichloroethene	(70)	4 J	1 J	2.00	1 J	4 U	1 U	4 U	4 U	4 U	4 U	4 U
Tetrachioroethene	(5)	1 J	3 J	4.00	3 J	0.6 J	0.7 J	0.6 J	1_U	1 U	1 U	1 U
Trichloroethene	(5)	1 J	2 J	3.00	2 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U

Highlighting indicates compound exceeded the Screening Level.

- U Constituent not detected; method detection limit (MDL) of the analysis is reported.
- J Concentration reported is an estimated value.
- UJ The analyte was not detected above the reported sample quantitation limit.
- DL Designates sample was diluted.
- B Designates the constituent was detected in the method blank.
- --- Indicates compound not analyzed.

Evergreen Manor

Remedial Investigation Report Section: 7

Revision: 1

Date: 28 March 2001

Page: 17 of 20

### Table 7-4 **Summary of Detected Constituents in Groundwater** Residential Wells - VOCs Evergreen Manor, Roscoe, Illinois

Sample ID:		RW01-01	RW02-01	RW-03	RW-04	RW04-01	RW-05	RW-05DUP	RW-06	RW-07	RW-08,
Sample Number:		_		-		EABQ4			-		-
Sample Date:		5/31/00	5/31/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00
Address:											<b></b>
Laboratory:		ESAT	ESAT	ESAT	ESAT	MITKEM	ESAT	ESAT	ESAT	ESAT	ESAT
Units		ug/L	ug/L	ug/L	ug/L	ug/L.	ug/L	ug/L	ug/L	ug/L	ug/L
PARAMETER (Screen	ing Level)										
1,1,1-Trichloroethane	(200)	2 U	2 U	2 J	2 J	2.00	2 U	2 U	2 U	5 J	0.9 J
Acetone	(700)	2 U	2 U	0.80	2 U	5 U	2 U	0.60	2 U	2 U	2 U
Chloroform	(0.02)	3 U	3 U	3 U	3 U	1 U	3 U	3 U	3 U	3 U	0.9 J
Cis-1,2-Dichloroethene	(70)	4 U	4 U	4 U	1 J	2.00	4 U	4 U	4 U	4 U	4 U
Tetrachloroethene	(5)	1 U	1 U	1 U	2 J	2.00	1 U	1 U	1 U	1 U	1 U
Trichloroethene	(5)	1 U	1 U	1 U	5 J	6.00	1 U	1 U	1 U	0.7 J	1 U

Section: 7 Revision: 1

Date: 28 March 2001

Page: 19 of 20

## Table 7-4 Summary of Detected Constituents in Groundwater Residential Wells - VOCs Evergreen Manor, Roscoe, Illinois

Sample ID:	RW-18	RW-19	RW-19DUP	RW-20	RW-21	RW-22		
Sample Number:								
Sample Date:	6/6/00	6/6/00	6/6/00	6/6/00	6/6/00	6/6/00		
Address:								
Laboratory:		ESAT	ESAT	ESAT	ESAT	ESAT	ESAT	
Units		ug/L	ug/L	ug/L.	ug/L	ug/L	ug/L	
PARAMETER (Screeni	ing Level)							
1,1,1-Trichloroethane	(200)	2 U	1 J	0.6 J	2 U	2 U	2 U	
Acetone	(700)	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ	2 UJ	
Chloroform	(0.02)	3 U	3 U	3 U	3 U	3 U	3 U	
Cis-1,2-Dichloroethene	(70)	4 U	4 U	4 U	4 U	4 U	4 U	
Tetrachloroethene	(5)	1 U	0.9 J	0.9 J	1 U	1 U	1 U	
Trichloroethene	(5)	1 U	4 J	4 J	1 U	1 U	1 Ü	

Highlighting indicates compound exceeded the Screening Level.

- U Constituent not detected; method detection limit (MDL) of the analysis reported.
- J Concentration reported is an estimated value.
- UJ The analyte was not detected above the reported sample quantition limit.
- B Designates the constituent was detected in the method blank.

Section: 7 Revision: 1

Date: 28 March 2001

Page: 18 of 20

## Table 7-4 Summary of Detected Constituents in Groundwater Residential Wells - VOCs Evergreen Manor, Roscoe, Illinois

Sample ID:		RW-10	RW-11	RW-11DUP	RW11-01	RW11-01DP	RW-12	RW-13	RW-14	RW-15	RW-16
Sample Number:					EABQ5	EABQ6					
Sample Date:	<del></del>	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/6/00	6/6/00	6/6/00	: 6/6/00
Address:											
Laboratory:		ESAT	ESAT	ESAT	MITKEM	MITKEM	ESAT	ESAT	ESAT	ESAT	ESAT
Units		ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
PARAMETER (Screen	ing Level)										
1,1,1-Trichloroethane	(200)	2 U	1 J	1 J	1.00	1.00	2 U	2 U	2 U	2 U	2 U
Acetone	(700)	2 U	2 U	2 U	5 U	5 U	2 U	2 UJ	2 UJ	2 UJ	2 UJ
Chloroform	(0.02)	3 U	3 U	3 U	1 U	1 U	3 U	3 U	3 U	3 U	3 U
Cis-1,2-Dichloroethene	(70)	4 U	4 U	4 U	1 U	1 U	4 U	4 U	4 U	4 U	4 U
Tetrachloroethene	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	(5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U

Section: 7 Revision: 1

Date: 28 March 2001 Page: 6 of 20

## Table 7-1 Screening Levels for Detected Constituents Evergreen Manor, Roscoe, Illinois

#### **GROUNDWATER**

GROUNDWATER				
	IEPA	U.S. EPA		
Constituent	Tier I Groundwater	Maximum Contaminant	Screening Level	
Constituent	Remediation Objective	Level		
	(μ <b>g/L)</b>	(μ <b>g/L</b> )	(μ <b>g/L</b> )	
1,1,1-trichloroethane	200	200	200	
1,1,2-trichloro-1,2,2-trifluoromethane			-	
1,1-dichloroethane	700		700	
1,2-dichloroethane	5	5	5	
2-butanone				
Acetone	700		700	
Benzene	5	5	5	
Chloroform	0.02		0.02	
cis-1,2-dichloroethene	70	70	70	
Ethylbenzene	7.00	700	700	
Methylene chloride	5	5	5	
Tetrachloroethene	5	5	5	
Toluene	1,000	1,000	1,000	
Trichloroethene	5	5	5	
Xylenes	10,000	10,000	10,000	

#### SEDIMENT

Constituent	IEPA Tier I Soil Remediation <b>Objecti</b> ve (µg <b>/kg</b> )	Region IX Risk Based Concentrations ** (µg/kg)	Screening Level (μg/kg)
Benzene	800 *	1,400	800
Chloroform	300 *	520	300
Methyl Acetate			
Toluene	650,000 *	520,000	520,000

- \* Remediation objective is based on the inhalation exposure route.
- \*\* Region IX RBCs are based on residential property use.

Revision: 1 ate: 28 March 2001

Date: 28 March 2001 Page: 7 of 20

### Table 7-2 Summary of Detected Consituents in Groundwater CPT - VOCs

### Evergreen Manor, Roscoe, Illinois

Sample ID:	CPT-01-01	CPT-01-02	CPT-01-03	CPT-01-04	CPT-01-05	CPT-01-06	CPT-01-07	CPT-01-08	CPT02-01	
Sample Number:							-	-	EABX3	
Sample Date:	5/25/00	5/25/00	5/25/00	5/25/00	5/25/00	5/25/00	5/25/00	5/25/00	6/2/00	
Laboratory:	ESAT	MitKem								
Sample Screen Depth (ft below ground	: 99	89	79	69	59	49	39	29	28	
Units:	ug/L	ug/L								
Parameter (Screening Level)										
Acetone (700)	2 UJ	15 B	2 UJ	5 U						
Methylene Chloride (5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	2 U	
1,1-Dichloroethane (700)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	
2-Butanone ()	11 U	5 U								
Cis-1,2-Dichloroethene (70)	4 U	2 J	1 J	1 J	1 J	1 J	4 U	4 U	1 U	
1,1,1-Trichloroethane (200)	2 U	1 J	1 J	1 J	1 J	2 U	1 J	2 U	1 U	
Benzene (5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	
Trichloroethene (5)	2 J	3 J	4 J	4 J	3 J	4 J	3 J	2 J	1 Ü	
Toluene (1000)	0.7 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.5 J	
Tetrachloroethene (5)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	
Ethylbenzene (700)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	
m-&/or p-Xylene (10,000)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U		
o-xylene (10,000)	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U		
Xylenes (total) (10,000)									0.6 J	

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RFW036-2A-AHVH

> Section: 7 Revision: 1

Date: 28 March 2001

Page: 20 of 20

# Table 7-5 Summary of Detected Constituents in Sediment VOCs Evergreen Manor, Roscoe, Illinois

Sample ID:	SD01-01	SD01-01DUP	SD02-01	SD03-01	SD04-01	SD04-01	SD04-01DP	SD05-01	SD06-01	
Sample Number:				_		EABQ9	EABR1	· ·		
Sample Date:	5/24/00	5/24/00	5/23/00	5/23/00	5/24/00	5/24/00	5/24/00	5/24/00	5/23/00	Screening
Laboratory:	ESAT	ESAT	ESAT	ESAT	ESAT	DATACHEM	DATACHEM	ESAT	ESAT	Level
Units	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg
PARAMETER										
Benzene	5 U	5 U	5 U	5 U	5 U	12 U	2 J	5 U	5 U	800
Chloroform	8 J	50 UJ	50 UJ	50 UJ	50 UJ	12 U	12 U	50 UJ	50 UJ	300
Methyl Acetate	5 J	15 U	9 J	15 U	15 U	12 U	12 U	9 J	15 U	
Toluene	5 U	5 U	5 U	5 U	5 U	12 U	0.7 J	5 U	5 U	650,000

- --- Screening Level is not available for this constituent.
- U Constituent not detected; method detection limit (MDL) of the analysis reported.
- J Concentration reported is an estimated value.
- UJ The analyte was not detected above the reported sample quantition limit.

Section: 8 Revision: 1

Date: 28 March 2001

Page: 1 of 21

**SECTION 8** 

**CONTAMINANT FATE AND TRANSPORT** 

The fate and transport of the contaminants present at the Evergreen Manor site, and the primary

geochemical factors influencing their concentrations and behavior, are discussed in this section.

These factors include dispersion, dilution, adsorption, oxidation, and geochemical behavior.

Physical characteristics of the site and the nature and extent of contamination, which have a

substantial influence on the factors affecting the fate and transport of contaminants, are also

addressed in this section. This section concludes with the results of a contaminant transport model

used to estimate the time it will take to achieve screening levels at the site.

8.1 CONTAMINANTS OF POTENTIAL CONCERN

The COPCs at the Evergreen Manor site were identified based on the extent of contamination at the

site and the contaminants' potential to migrate. The COPCs at the site include three VOCs in

groundwater, as described in Section 7 and the following subsections.

Groundwater

Chemical constituents exceeding screening levels in groundwater include chloroform,

tetrachloroethene, and trichloroethene. Chloroform was detected at a concentration considerably

above its screening level, and tetrachloroethene and trichloroethene were detected only slightly

above their screening levels; therefore, these three VOCs are considered to be COCs in the

groundwater at the Evergreen Manor site.

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RFW036-2A-AHVH

Section: 8 Revisión: 1

Date: 28 March 2001

Page: 2 of 21

8.2 FATE OF CONTAMINANTS OF CONCERN

In groundwater, chloroform, tetrachloroethene, and trichloroethene were detected above their

screening levels. These compounds are classified as volatile chlorinated hydrocarbons, and are

moderately to readily water soluble.

Chloroform has an adsorption rate to soil that should be insignificant at the Evergreen Manor site.

Therefore, chloroform should be highly mobile in the groundwater environment. Chloroform does

not readily biodegrade in groundwater and may remain in the dissolved phase for extended periods

of time. Bioconcentration is not expected to be significant. Biodegradation products include

methylene chloride (Howard, 1990).

Tetrachloroethene has a moderate adsorption rate to soil that creates a low to medium mobility in

the groundwater environment. Biodegradation does not occur under aerobic conditions and is slow

under anaerobic conditions if the microbes have been acclimated. Bioconcentration is not expected

to be significant in aquatic organisms. Biodegradation products include trichloroethene, cis- and

trans-1,2-dichloroethene, methylene chloride, chloroethene, and vinyl chloride (Howard, 1990).

Trichloroethene has an adsorption rate to soil that should not be significant. Therefore,

trichloroethene is expected to be highly mobile in the groundwater environment. Biodegradation

occurs in water under most conditions. Bioconcentration is moderate in aquatic organisms.

cis-1,2-Dichloroethene is the primary biodegradation product (Howard, 1990).

Physical and chemical properties of these organic compounds are presented in greater detail in Table

8-1.

Section: 8 Revision: 1

Date: 28 March 2001

Page: 3 of 21

8.3 CONTAMINANT MIGRATION PATHWAY

After a chemical is released into the environment, it may be transported (i.e., advective transport),

physically transformed (i.e., volatilized), chemically transformed (i.e., via oxidation/reduction),

biologically transformed (i.e., biodegradation), or bioaccumulated in one or more media.

Contaminant migration pathway analysis identifies other (non-source) environmental media and off-

site areas potentially affected by contaminant migration.

The following addresses the possible migration pathways of contaminants at the Evergreen Manor

site, their potential to be transported to other environmental media, and their potential to migrate off-

site. Table 8-2 presents a summary of the migration characteristics of contaminants at the site.

Groundwater Dispersion and Biodegradation

Most of the contaminant transport at the site is achieved through advective transport (i.e., through

groundwater movement). However, because of the low concentrations present at the site, it is likely

that the concentrations are also being reduced through dispersion and/or biodegradation. Dispersion

can be thought of as the spread of contaminants that occurs in addition to being transported by

advection. Biodegradation is the chemical break down of one constituent into another, or into a

more elemental form. The result of both of these mechanisms is to reduce the concentration of the

constituents in groundwater.

Table 8-3 presents detected constituent concentrations from the HRS package and this RI. Average

concentrations are also presented and show the decline in groundwater VOC levels between the HRS

scoring period and the current RI results. Table 8-4 directly compares residential well data from the

HRS package and this RI for locations sampled during both sampling events. The two comparable

locations show that 1,1,1-TCA and TCE concentrations have declined, while the trend is unknown

for cis-1,2-DCE and PCE, since data were not presented for these locations in the HRS package.

Table 8-5 compares monitoring well data from the HRS package and this RI for wells sampled

during both events. In general the contaminant concentrations declined, or in some instances

CHLAN01\WP\WO\RAC\036\29672S-8.WPD

RFW036-2A-AHVH

124

Section: 8 Revision: 1

Date: 28 March 2001

Page: 4 of 21

remained approximately the same. The trend is not known for cis-1,2-DCE and PCE concentrations

at MW-104S, because the HRS package did not present data for these constituents.

The effectiveness of biodegradation was evaluated by following a procedure found in the Technical

Protocol for Evaluating Natural Attenuation of Chlorinated Solvents in Ground Water (U.S. EPA,

1998c). By applying a score to the various water quality parameters and constituent concentrations,

at each sampling point, it was possible to judge whether or not anaerobic biodegradation (reductive

dechlorination) could be occurring. The highest cumulative score achieved was three points, which

was interpreted as, "Inadequate evidence for anaerobic biodegradation of chlorinated organics," (U.S.

EPA, 1998c). Strong evidence of biodegradation can be achieved with a score of 15 or more points.

Thus, biodegradation is not expected to be a considerable factor to concentration reduction of COCs,

when compared to the effect of dispersion. This does not imply, however, that biodegradation is not

occurring.

Groundwater to Air Pathway

Based on groundwater depths and chemical contaminants present, it is unlikely that migration of

chemical contaminants from groundwater to air would occur. Volatilization of VOCs in

groundwater is possible; however, this is not likely to be a significant process and is subsequently

of minimal concern.

Therefore, the migration of groundwater contaminants via the groundwater to air pathway is not a

pathway of potential concern.

Groundwater to Surface Water Pathway

Contaminants in groundwater may migrate to the Rock River. Because of the difference in elevation

between the water table and Dry Creek, in the vicinity of the site, it is not expected that contaminated

groundwater would migrate to Dry Creek. Based on the groundwater gradient shown in Figure 6-5.

and the elevation of the Rock River south of the site (approximately 700 ft MSL), the expectation

CHLAN01\WP\WO\RAC\036\29672S-8.WPD

RFW036-2A-AHVH

Section: 8 Revision: 1

Date: 28 March 2001

Page: 5 of 21

is that groundwater discharges to the Rock River. Migration of contaminants could occur where

groundwater discharges to surface water.

The results of the surface water sampling indicated that VOCs were not detected in the Rock River.

This is most likely due to dilution that occurs when a relatively small volume of groundwater is

discharged to the Rock River and is mixed with a relatively large volume of surface water.

Based on the relatively low COC concentrations detected at the site, and the large amount of dilution

occurring, the groundwater to surface water migration pathway does not appear to be a concern at

the Evergreen Manor site.

8.4 TRANSPORT MODELING

A modeling approach was taken to estimate the time for contaminants to decline to below screening

levels. A simple groundwater model was used to simulate the transport of contaminants through the

saturated subsurface. Based on the available data, an analytical model approach was determined to

be applicable. The BIOSCREEN Natural Attenuation Decision Support System (Newell, 1996) was

the model used to simulate contaminant transport.

8.4.1 BIOSCREEN

BIOSCREEN was written to support natural attenuation of hydrocarbons at petroleum sites,

however, the transport code is equally applicable for other dissolved contaminants. The model takes

into consideration advection, dispersion, adsorption, and biodegradation; however, since

biodegradation could not be proved to be occurring at the Evergreen Manor site, based on RI

analytical data, it was not incorporated into the model.

BIOSCREEN models a single contaminant originating from a source area with a known contaminant

mass. The model uses a half-life approach to reducing the contaminant mass at the source. With

a small source mass input, the model can be used to approximate a short term or nearly instantaneous

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24

Section: 8 Revision: 1

Date: 28 March 2001

Page: 6 of 21

contaminant release. Although unknown, it is presumed that the release at the site can be modeled

as short term or nearly instantaneous release.

The limitations of the BIOSCREEN model are that it assumes simple groundwater flow conditions,

and only approximates more complicated processes. The sand and gravel aquifer underlying the

Evergreen Manor site is assumed to be fairly homogeneous, and can be modeled as one continuous

flow system. The distribution of chemical data at the Evergreen Manor site is more complex, and

BIOSCREEN was used to provide approximations of contaminant concentrations.

BIOSCREEN can estimate concentration distributions either along the axis of a plume, or across the

modeled area. To simplify the modeling approach, calibration data were assumed to be located

along the axis of the plume, and only the output of concentration distributions along the axis of the

plume were evaluated.

8.4.2 Input and Assumptions

Table 8-6 presents the input parameters used in each of the four models created. Since these

parameters vary for each contaminant, a separate model was created for each one. The contaminants

include chloroform, 1,1,1-TCA, TCE, and PCE.

Model Dimensions - Assuming that the contaminant source area is located in the area north of

Rockton Road, and east of IL 251, the length of the plume (L<sub>n</sub>) was set to be 13,000 feet. This is the

straight line length to the presumed discharge area at the Rock River. A width of 2,500 feet was

used, which is approximately twice the presumed width of the actual plume.

Source Concentration and Source Mass - Since a source has not been identified at the site, the

source concentration and source mass were adjusted during modeling to fit the calibration data.

These values were altered for each compound.

Hydraulic Conductivity - The value provided in Section 6 of this report is 3.8 X 10<sup>-2</sup> cm/sec. This

value is based on pressure tests, which are similar to slug tests in that they only approximate the

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Section: 8 Revision: 1

Date: 28 March 2001

Page: 7 of 21

hydraulic conductivity of the aquifer volume in close proximity to the test well. It is possible that the volume of aquifer close to the borehole has been disturbed during well installation (Kruseman, 1990), or has differing hydraulic properties, and the resulting hydraulic conductivity could be underestimated. The hydraulic conductivity used in modeling was  $2.2 \times 10^{-1}$  cm/sec, which best fit the modeled concentrations to the available data and is a value within the range of hydraulic conductivities for the types of geologic materials found at the site.

Hydraulic Gradient - The value used, 0.0015 ft/ft, is based on the groundwater elevation data presented in Section 6.

Porosity - The value used, 30%, is a typical porosity for sand and gravel mixtures (Fetter, 1994).

**Dispersion** - For chloroform, 1,1,1-TCA, and TCE, longitudinal dispersivity (Alpha X) was set to 59.9 feet, and was calculated with the Xu and Eckstein (1995) equation:

$$AlphaX = 3.28 \cdot 0.83 \left[ \log_{10} \left( \frac{L_p}{3.28} \right) \right]^{2.414}$$

This equation is based on the length of the plume, which equals 13,000 feet. The longitudinal dispersivity was set to 100 feet for the PCE model. Transverse dispersivity was set to one-tenth of the longitudinal dispersivity. Vertical dispersivity was anticipated to be negligible compared to longitudinal and transverse dispersion.

Retardation Factor - This was calculated using a soil bulk density ( $\rho_b$ ) of 1.8 kg/L (IAC, 1997), a contaminant specific partition coefficient ( $K_{oc}$ ), a fraction of organic carbon ( $f_{oc}$ ) of either 0.06% or 0.2% (U.S. EPA, 1998b), and a porosity (n) of 30% in the following equations:

$$K_d = K_{oc} \cdot f_{oc} \qquad R_f = 1 + \frac{K_d \cdot \rho_b}{n}$$

Section: 8 Revision: 1

Date: 28 March 2001

Page: 8 of 21

where  $K_d$  is the contaminant-specific distribution coefficient. The partition coefficient values were

obtained from Table 8-1 and from Groundwater Chemicals Desk Reference (Montgomery, 1989).

Retardation factors of 1.5, 2.5, 2.1 and 2.0 were used for chloroform, 1,1,1-TCA, TCE, and PCE,

respectively.

8.4.3 Calibration Data

The four contaminants modeled with BIOSCREEN included chloroform, 1,1,1-TCA, TCE, and PCE.

Although concentrations of 1,1,1-TCA did not exceed screening levels, this contaminant was used

in order to calibrate the model. Table 8-7 presents the data to which the models were calibrated.

This table presents the concentrations from the HRS package and this RI, as well as the approximate

distance from the source area. The HRS package data were collected about 5 to 6 years prior to the

RI data. This time interval was also used to calibrate the models.

8.4.4 1,1,1-TCA and TCE

The 1,1,1- TCA and TCE models were calibrated to the data presented in Table 8-7 by altering the

hydraulic conductivity, the source concentration, and the source mass. Several attempts were made

to match the calibration data using a hydraulic conductivity value of 3.8 X 10<sup>-2</sup> cm/sec, as presented

in section 6-2, however, the calibration data could not be matched. As discussed previously, the

hydraulic conductivity could have been underestimated. Therefore, the hydraulic conductivity was

increased until the modeled contaminant distribution matched the calibration data, yet still resulted

in using an acceptable value with regard to geologic conditions.

The source concentration and the source mass were adjusted in order to approximate the actual

concentrations from the HRS package and this RI. Although an actual source concentration or mass

are not known, values were chosen that fitted the available data. The resulting concentration

distributions are presented in Appendix F.

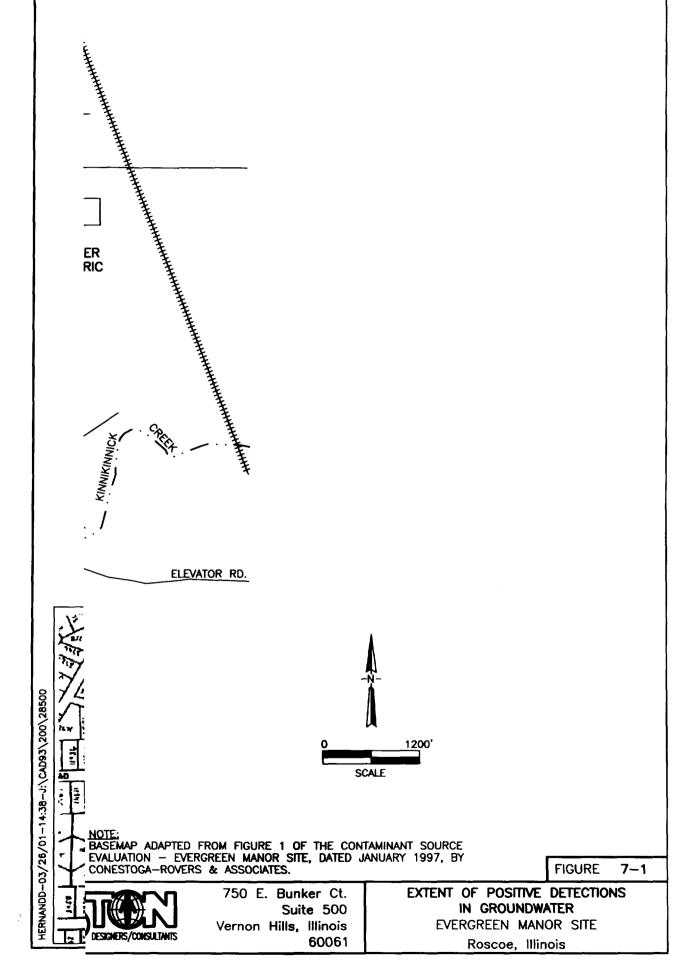
The plots for 1,1,1-TCA indicate that the HRS package data and the RI data can be matched at 24

and 30 years respectively. These times do not represent the actual time since a release occurred, but

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Section: 8 Revision: 1

Date: 28 March 2001

Page: 9 of 21

rather the time it took to match the calibration data. Thus, the 24 year output matches the HRS data,

and the 30 year output matches the RI data. Similarly, the TCE concentrations were matched to the

HRS package and RI data at model output at 20 and 26 years.

Simulations of TCE transport were run beyond the RI time frame to estimate when concentrations

would decline below the screening level of 5 ug/L. This result was achieved from the 32 year

simulation, or 6 years after the RI. Thus, in about 2006 TCE concentrations at the site are predicted

to be below the screening level.

8.4.4 Chloroform

As shown on Table 8-7, only one data point exists for chloroform. Based on the calibrations

performed for 1,1,1-TCA and TCE, a model for chloroform was created which matched the

concentration at 11943 Wagon Lane during the RI. Only the adsorption, source concentration, and

source mass values were adjusted to model chloroform transport. The result, presented in

Appendix F, shows that after a simulated time of 15 years, the RI datum is matched.

The simulation was run beyond the 15 years to estimate when the chloroform concentration would

decline below the screening level of 0.02 ug/L. This result was achieved with the 18 year

simulation, or 3 years after the RI. Thus, chloroform concentrations at the site are predicted to be

below the screening level in 2003.

8.4.5 PCE

The PCE model was created by matching three calibration data points from the RI. Dispersion,

adsorption, source concentration, and source mass values were adjusted to model PCE transport.

The model output at 15 years corresponds to the data collected during the RI. This result is

presented as part of Appendix F.

The simulation was run beyond the 15 years to estimate when PCE concentrations would fall below

the screening level, 5 ug/L. The 30 year simulation was found to meet this goal. Thus, about 15

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Evergreen Manor Remedial Investigation Report Section: 8 Revision: 1 Date: 28 March 2001

Page: 10 of 21

years after the RI, in 2015, PCE concentrations at the site are predicted to be below the screening level.

Evergreen Manor

Remedial Investigation Report

Section: 8 Revision: 1

Date: 28 March 2001 Page: 11 of 21

Table 8-1

Physical and Chemical Properties of Constituents of Potential Concern vergreen Manor Roscoe, Illinois

Constituent	Hazard Class	Physical Form	Physical Description	General Chemical Class	Molecular Weight (g/mole)	Log K <sub>m</sub> *	Log K <sub>us</sub> t	Density <sup>e</sup> (unitless)	Boiling Point (°C)	Viscosity	Water Solubility (mg/L)	Henry's Law Constant (stm-m³/mol)	Vapor Pressure (mm Hg)	Flash Point (°C)
Volatile Organic Cos														
Chloroform	Poison	Liquid	Clear, water-white volatile liquid	Solvent	119.38	1.64	1.94	1.4861	61.7	0.542 @ 25°C	8,380 @ 25°C	0.0032	198 @ 25°C	Noncom- bustible
Trichloroethene	Poison	Liquid	Clear, colorless watery liquid w/ chloroform-like odor.	Solvent	131.39	1.98	2.59	1.463	87.0	NA	1,285 @ 25°C	0.010	73.3 @ 25°C	32.2
Tetrachloro-ethene	Poison	Liquid	Colorless liquid w/ sweet etheral odor	Solvent	165.83	2.43	2.53	1.6226	121.2	NA	345 @ 25°C	0.0142	20 @ 25°C	Not Flammable

#### Notes:

NA - Not available.

N/A - Not applicable.

CHLAN01\WP\WO\RAC\036\29672T8-1.WPD

RFW036-2A-AHVH

<sup>\*</sup> Organic carbon partition coefficient.

<sup>&</sup>lt;sup>b</sup> Octanol-water partition coefficient.

<sup>&</sup>lt;sup>c</sup> Density of the compound at 20°C in relation to water at 4°C, unless specified.

Section: 8 Revision: 1

Date: 28 March 2001 Page: 12 of 21

Table 8-2 Migration Characteristics of Constituents of Potential Concern Evergreen Manor Roscoe, Illinois

Constituent	Sorption	Biodegradability/Bioc Biotransform		Photodegradation Rates	Hydrolysis Rates	Chemical Transformations
Volatile Organ	ic Compounds (VOCs)					
Chloroform	Adsorbs most strongly to peat moss, but not at all to sand. Should not be adsorbed readily at the Evergreen Manor site.	Very slow biodegradation anaerobic conditions have when microbes have acclin chemical / there is little to bioconcentration potential biodegradation products in chloride.	en reported ated to the	Photodegradation is not a significant loss process in aquatic systems, but is slow in atmospheric conditions with hydroxly radicals, with a half-life of 80 days.	Has a negligible rate of hydrolysis.	Degradation (loss of a chloride ion) will result in methylene chloride formation. Minor sources of chloroform release include, but are not limited to, the decomposition of trichloroethylene.
Trichloro- ethene	Very weak adsorption to most soils.	Slow biodegradation in wa conditions / moderate biocaquatic organisms / biodeg include cis- and trans-1,2-a and vinyl chloride.	centration in dation products	Direct photolysis does not occur; reaction occurs with hydroxyl radicals in the atmosphere, with a half-life of 5 days.	Hydrolysis does not occur under normal conditions.	Reaction with hydroxyl radicals in the atmosphere produces phosgene, dichloroacetyl chloride, and formyl chloride.
Tetrachloro- ethene	Low to medium mobility in soil is expected; therefore, adsorption should be moderate.	There is no evidence for bi under aerobic conditions, biodegradation occurs undeconditions if the microbes acclimated / bioconcentrati expected to be significant i organisms / biodegradation trichloroethene, cis- and tridichloroethene, methylene chloroethene, and vinyl chl	t slow anaerobic ive been n is not aquatic products include s- 1,2- hloride,	Vapor-phase reaction with photochemically produced hydroxyl radicals occurs with a half-life of 2 months.	Not expected to significantly hydrolyze in soil or water.	Slow biodegradation under anaerobic conditions when the organisms have been acclimated yields trichloroethene. Traces of dichloroethylene isomers and vinyl chloride were also found.

Table 8-3
Comparison of Previous Groundwater Data with RI Data
Evergreen Manor, Roscoe, Illinois

Well ID	1,1,1-TCA	TCE	1,2-DCE	PCE
Residenti	al Well Sample	Results Suppo	rting HRS Score	e (1993-1995)
G 103	19	31	n/a	n/a
G 104	15	23	n/a	n/a
G 105	13	20	n/a	n/a
G 106	14	23	n/a	n/a
G 107	24	35	n/a	n/a
G 108	15	20	n/a	n/a
G 109	10	17	n/a	n/a
G 110	10	18	n/a	n/a
G 112	12	23	n/a	n/a
G 113	23	38	n/a	n/a
G 114	20	36	n/a	n/a
G 115	15	27	n/a	n/a
G 116	18	27	n/a	n/a
G 117	12	24	n/a	n/a
G 118	n/a	19	n/a	n/a
G 119	29	19	n/a	n/a
G 120	14	24	n/a	n/a
G 121	10	18	n/a	n/a
G 122	17	25	n/a	n/a
G 123	22	23	n/a	n/a
G 124	17	30	n/a	n/a
G 125	n/a	19	n/a	n/a
G 129	14	25	n/a	n/a
	, , , , i	1,		·
G 131	22	20	n/a	n/a
G 132	12	18	n/a	n/a
G 134	16	25	n/a	n/a
G 135	21	29	n/a	n/a
G 136	24	29	n/a	n/a
G 137	20	27	n/a	n/a
G 138	12	23	n/a	n/a
G 139	n/a	11	n/a	n/a
G 141	10	18	n/a	n/a
G 142	33	40	n/a	n/a
G 143	37	34	n/a	n/a
G 144	34	35	n/a	n/a
G 146	21	28	n/a	n/a
G 147	11	22	n/a	n/a
G 148	22	31	n/a	n/a
G 149	13	22	n/a	n/a
G 152	18	6	2	2
G 153	21	15	5	4
G 154	18	22	8	5
G 155	21	10	3	2

Table 8-3
Comparison of Previous Groundwater Data with RI Data
Evergreen Manor, Roscoe, Illinois

Well ID	1,1,1-TCA	TCE	1,2-DCE	PCE
	I	Results Suppo	rting HRS Scor	e (1993-1995)
G 156	21	6	2	2
G 157	17	6	2	2
G 167	22	12	3	2
G 168	16	13	3	n/a
G 170	30	6	2	n/a
G 184	17	10	2	4
G 188	11	2.1	n/a	n/a
G 190	14	27	5.3	n/a
G 192	1.6	6	n/a	n/a
G 196	28	5.7	n/a	n/a
G 202	19	7.3	1	n/a
G 203	20	8.2	1.2	n/a
G 206	34	15	n/a	n/a
G 212	21	n/a	n/a	n/a
G 219	5.1	11	1.4	n/a
G 229	19	20	2.8	n/a
G 241	26	n/a	n/a	n/a
G 247	20	31	4.3	n/a
G 248	25	38	4.7	n/a
G 251	n/a	n/a	3.8	n/a
G 257	n/a	5.4	n/a	n/a
G 260	12	12	1.2	n/a
G 268	18	13	1.9	n/a
U 47U	14	د ـ	7	104
G 276	25	23	3.1	n/a
G 283	16	11	1.6	n/a
G 290	17	22	3.1	n/a
G 293	14	12	1.4	n/a
G 296	18	19	2.5	n/a
G 304	19	25	n/a	n/a
G 316	16	31	4.4	n/a
G 317	29	24	2.8	n/a
G 318	9.7	17	2.2	n/a
G 320	11	5.3	0.6	n/a
G 322	3.6	11	2.3	n/a
G 338	16	23	6.4	4
G 357	14	12	1.6	n/a
G 358	21	28	3.4	n/a
G 360	16	28	4.1	n/a
G 362	26	27	4	n/a
G 364	16	11	1.8	n/a
G 365	16	17	2.7	n/a
Average:	18	20	3	3

Table 8-3
Comparison of Previous Groundwater Data with RI Data
Evergreen Manor, Roscoe, Illinois

Well ID	1,1,1-TCA	TCE	1,2-DCE	PCE						
	Residential Well Sample Results from RI (2000)									
RW 3	2	ND	ND	ND						
RW 4	2	5	1	2						
RW 4	2	6	2	2						
RW 7	5	0.7	ND	ND						
RW 8	0.9	ND	ND	ND						
RW 11	1	ND	ND	ND						
RW 11	1	ND	ND	ND						
RW 11	1	ND	ND	ND						
RW 11	1	ND	ND	ND						
RW 19	1	4	ND	0.9						
RW 19	0.6	4	ND	0.9						
Average:	1.6	3.9	1.5	1.5						

Monitorin	g Well Sample	Results Suppo	rting HRS Scor	e (1994-1995)
G 101D	n/a	3	n/a	n/a
G 103S	3	n/a	n/a	40
G 104S	12	0.9	n/a	n/a
G 105D	8.9	15	5.7	3.2
G 105D	9	15	5	4
G 106S	1	3	n/a	n/a
G 107D	8	n/a	n/a	11
G 108D	7	n/a	n/a	3
G 109D	8	3	6	7
G 110S	4	2	n/a	n/a
G 113	n/a	n/a	n/a	2
G 114	3	n/a	п/а	n/a
Average:	6.4	6.0	5.6	10
	Monitoring W	<mark>ell Sample R</mark> esi	ults from RI (20	00)
MW 101D	ND	2	ND	ND
MW 103D	3	ND	ND	0.5
MW 103S	ND	ND	ND	9
MW 104D	1	ND	ND	ND
MW 104D	1	ND	ND	ND
MW 104S	2	ND	ND	ND
MW 105D	2	2	1	3
MW 105D	3	3	2	4
MW 105S	2	2	1	3
MW 108D	ND	ND	ND	0.6
MW 108D	ND	ND	ND	0.7
MW 108D	ND	ND	ND	0.6
MW 112	2	ND	ND	ND
Average:	2.0	2.3	1.3	2.7

n/a - Data not available, not analyzed, or compound not detected.

ND - Compound not detectd

Section: 8 Revision: 1 Date: 28 March 2001 Page: 16 of 21

## Table 8-4 Comparison of Sele t Detected Constituents in Groundwater R: sidential Wells - VOCs

Evergeen Manor, Roscoe, Illinois

THE RESERVE OF THE PARTY OF THE							agreement of the state of the state of the state of		**** ***				
Sampling Event:	RI	RI	RI	HRS	RI	RI	RI	RI	RI	RI	RI	RI	HRS
Sample ID:	RW-03	RW-04	RW04-01	G109	RW-(	RW-08	RW-11	RW-11DUP	RW11-01	RW11-01DP	RW-19	RW-19DUP	G110
Sample Number:	_	_	EABQ4	_		I –	_		EABQS	EABQ6	_	_	-
Sample Date:	6/5/00	6/5/00	6/5/00	11/9/93	6/5/0	6/5/00	6/5/00	6/5/00	6/5/00	6/5/00	6/6/00	6/6/00	11/9/93
Address:	11990 Wagon La.	11990 Blue Spruce	11990 Blue Spruce	11990 Blue Spruce	12031 W 20m La. C	11943 Wagon La.	4234 Valorio Rd.	4234 Valorio Rd.	4234 Valorio R4.	4234 Valerie Rd.	11974 Blue Spruce Dr.	11974 Blue Spruce Dr.	11974 Blue Spruce Dr.
Laboratory:	ESAT	ESAT	MITKEM	Unknown	ESA	ESAT	ESAT	ESAT	MITKEM	MITKEM	ESAT	ESAT	Unknown
Units	ug/L	ug/L	ug/L	ug/L	ug/1	ug/L	ug/L	•g/L	wg/L	wg/L	<b>ug/</b> L	ug/L	ug/L
PARAMETER													
1,1,1-Trichioreethane	2 J	2 J	2.00	10	5 )	0.9 J	1 J	1 J	1.00	1.00	1 J	0.6 J	10
Cis-1,2-Dichloroethese	ND	1 J	2.00	n/a	NL	ND	ND	ND	ND	ND	ND	ND	n/a
Tetrachlereethene	ND	2 J	2.00	n/a	NI	ND	ND	ND	ND	ND	0.9 J	0.9 J	n/a
Trichiereethene	ND	5 J	6.00	17	0.7	ND	ND	ND	ND	ND	4 J	4 J	18

#### Notes and Qualifler Flags

- ND Constituent not detected above method detection limit.
- J Concentration reported is an estimated value.
- n/a Constituent not presented in HRS package; constituent is presumed either not analyzed, or . . alyais resulted in a non-detect.

Section: 8

Date: 28 March 2001 Page: 17 of 21

Revision: 1

### **Comparison of Select Detected Constituents in Groundwater**

Moi itoring Wells - VOCs Evergre in Manor, Roscoe, Illinois

Table 8-5

Sampling Event:	RI RI	HRS	RI	RI	RI_	RI	RI_	RI	HRS
Sample ID:	MW101D	G101D	MW1018	MW102D	MW102S	MW103D	MW103D-DL	MW103S	G103S
Sample Number:		_							
Sample Date:	5/30/00	2/23/95	5/30/00	6/1/00	6/1/00	5/31/00	5/31/00	5/31/00	2/21/95
Laboratory:	ESAT		ESAT	ESAT	ESAT	ESAT	ESAT	ESAT	
Units	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
PARAMETER									
1,1,1-Trichloroethane	ND	n/a	ND	ND	ND	3 J	ND	ND	3
Cis-1,2-Dichloroethene	ND	n/a	ND	ND	ND	ND	ND	ND	n/a
Tetrachloroethene	ND	n/a	ND	ND	ND	0.5 J	ND	9 J	40
Trichloroethene	2 J	3	ND	ND	ND	ND	ND	ND	n/a

#### Notes and Qualifler Flags

- ND Constituent not detected above the method detection limit.
- J Concentration reported is an estimated value.
- DL Designates sample was diluted.
- n/a Constituent not presented in HRS package; constituent is presumed either not a dyzed, or analysis resulted in a non-detect.

Evergreen Manor Remedial Investigation Report Section: 8 Revision: 1 Date: 28 March 2001

Page: 18 of 21

# Table 8-5 Comparison of Select Detected Constituents in Groundwater Monitoring Wells - VOCs Evergre :n Manor, Roscoe, Illinois

Sampling Event:	RI	RI	RI	HRS	RI	RI	HRS	HRS	RI
Sample ID:	MW104D	MW104D-DUP	MW104	G104S	MW105D	MW105D-01	G105D	G105D	MW105S
Sample Number:						EABX1			
Sample Date:	6/2/00	6/2/00	6/2/00	2/22/95	6/2/00	6/2/00	3/23/94	2/22/95	6/2/00
Laboratory:	ESAT	ESAT	ESAT		ESAT	MitKem	_		ESAT
Units	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
PARAMETER									
1,1,1-Trichloroethane	1 J	1 J	2 J	12	2 J	3	8.9	9	2 J
Cis-1,2-Dichloroethene	ND	ND	4 J	n/a	1 J	2	5.7	5	1 J
Tetrachloroethene	ND	ND	1 J	n/a	3 J	4	3.2	4	3 J
Trichloroethene	ND	ND	1 J	0.9	2 J	3	15	15	2 J

#### Notes and Qualifler Flags

- ND Constituent not detected above the method detection limit.
- J Concentration reported is an estimated value.
- DL Designates sample was diluted.
- n/a Constituent not presented in HRS package; constituent is presumed either not nalyzed, or analysis resulted in a non-detect.

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> Section: 8 Revision: 1

Date: 28 March 2001 Page: 19 of 21

# Table 8-5 Comparison of Select Detected Constituents in Groundwater Mon toring Wells - VOCs Evergree 1 Manor, Roscoe, Illinois

Sampling Event:	RI	RI	RI	HRS	RI	RI	RI	HRS	RI
Sample ID:	MW108D	MW108D-01	MW108D-DUP	G108D	MW108S	MW110D	MW110S	G110S	MW112
Sample Number:		EABW8	-	_		<u> </u>			_
Sample Date:	6/1/00	6/1/00	6/1/90	2/21/95	6/1/00	6/1/00	6/1/00	2/23/95	6/2/00
Laboratory:	ESAT	MitKem	ESAT		ESAT	ESAT	ESAT		ESAT
Units	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
PARAMETER									
1,1,1-Trichloroethane	ND	ND	ND	7	ND	ND	ND	4	2 J
Cis-1,2-Dichloroethene	ND	ND	ND	n/a	ND	ND	ND	n/a	ND
Tetrachloroethene	0.6 J	0.7 J	0.6 J	3	ND	ND	ND	n/a	ND
Trichloroethene	ND	ND	ND	17/2	ND	ND	ND	2	ND

#### Notes and Qualifler Flags

- ND Constituent not detected above the method detection limit.
- J Concentration reported is an estimated value.
- DL Designates sample was diluted.
- n/a Constituent not presented in HRS package; constituent is presumed either not analy. d, or analysis resulted in a non-detect.

Section: 8 Revision: 1

Date: 28 March 2001 Page: 20 of 21

# Table 8-6 BIOSCREEN Model Inputs and Results Evergreen Manor, Roscoe, Illinois

Compound Modeled	Chloroform	1,1,1-TCA	TCE	PCE
Model Dimensions				
Length (ft)	13,000	13,000	13,000	13,000
Width (ft)	2,500	2,500	2,500	2,500
Source				
Width (ft)	500	500	500	500
Thickness (ft)	40	40	40	40
Concentration (mg/L)	0.006	0.045	0.06	0.022
Mass (kg)	0.15	35	50	60
Hydraulic Conductivity (cm/sec)	2.2E-01	2.2E-01	2.2E-01	2.2E-01
Hydraulic Gradient (ft/ft)	0.0015	0.0015	0.0015	0.0015
Porosity (dimensionless)	0.3	0.3	0.3	0.3
Seepage Velocity (ft/year)	1,138	1,138	1,138	1,138
Longitudinal Dispersion (ft)	59.9	59.9	59.9	100
Transverse Dispersion (ft)	6.0	6.0	6.0	10
Vertical Dispersion (ft)	1E-99	1E-99	1E-99	1E-99
Fraction Organic Carbon (mg/kg)	0.002	0.002	0.002	0.0006
Partition Coefficient (K <sub>oc</sub> ) (L/kg)	43.7	125.9	95.5	269.2
Distribution Coefficient (K <sub>d</sub> ) (dimensionless)	0.087	0.25	0.19	0.16
Soil Bulk Density (r <sub>b</sub> ) (kg/L)	1.8	1.8	1.8	1.8
Retardation Factor (dimensionless)	1.5	2.5	2.1	2.0
		Not		
Estimated Years to Reach Screening Level	) ]	Applicable	v	10

Evergreen Manor Remedial Investigation Report Section: 8 Section: 8 Removement

Date: 28 March 2001 Page: 21 of 21

# Table 8-7 BIOSCREEN Calibration Data Evergreen Manor, Roscoe, Illinois

Contaminant	Sample Location	Distance from Source (ft)	HRS Data (ug/L)	RI Data (ug/L)
Chloroform		11200		0.9 J
1,1,1-TCA	7	10,700	10	2
	-	8,850	9	3
TCE	•	10,700	17	6
	•	8,850	15	3
PCE		10,700		2
		2,700	_	9
		8,850		4

#### Notes:

J - Approximate concentration.

Section: 9 Revision: 1

Date: 28 March 2001

Page: 1 of 28

**SECTION 9** 

**HUMAN HEALTH RISK ASSESSMENT** 

9.1 INTRODUCTION

In accordance with the National Oil and Hazardous Substances Pollution Contingency Plan (NCP),

baseline human health and ecological risk assessments were prepared to evaluate the potential human

health and environmental impacts associated with the Evergreen Manor site in Roscoe, Winnebago

County, Illinois under a no-action alternative (i.e., in the absence of remedial [corrective] action).

Information and data collected as part of the RI/FS activities serves as the basis for these tasks.

Exposure to volatile organic compounds (VOCs) in groundwater is the primary focus of this risk

assessment.

9.1.1 Objectives

The objectives of the baseline human health risk assessment for the Evergreen Manor site are as

follows:

• Estimate potential risk to people contacting site-related chemicals of potential

concern(COPCs) under scenarios of current and plausible future land use.

Provide an analysis of risks and help determine the need for remedial actions at the

site.

Identify specific media and areas associated with unacceptable risk, if applicable.

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Date: 28 March 2001 Page: 2 of 28

Revision: 1

### 9.1.2 Risk Assessment Approach

The methodology used to assess the potential human health risks at the Evergreen Manor site draws upon the guidance set forth in the following documents:

- Risk Assessment Guidance for Superfund, Human Health Evaluation Manual, Part A (U.S. EPA, 1989).
- Risk Assessment Guidance for Superfund, Human Health Evaluation Manual, Part B (U.S. EPA, 1991a).
- Risk Assessment Guidance for Superfund, Human Health Evaluation Manual, Part D (U.S. EPA, 1998).

### 9.1.3 Risk Assessment Organization

The Human Health risk assessment is organized into the following components:

- <u>Hazard Identification</u>—Identification of major contaminants of concern based on a
  review of available information on the hazardous substances present at the site.
  COPCs were selected based on their intrinsic toxicological properties as part of a
  dose-response assessment.
- Exposure Assessment—Critical exposure pathways were identified and analyzed. The proximity of contaminants to exposure pathways and their potential to migrate into critical pathways was also assessed. Potential receptors were identified and characterized. The exposure assessment then identifies the magnitude of actual or potential human exposures, the frequency and duration of these exposures, and the exposure routes. The assessment also includes an evaluation of the likelihood of such exposures occurring, and provides the basis for the development of acceptable exposure levels. In developing the exposure assessment, reasonable maximum estimates of exposure for both current and future land used conditions were developed.
- <u>Toxicity Assessment</u>—Chemical-specific toxicity information is provided for the chemicals of potential concern.

Section: 9 Revision: 1

Date: 28 March 2001

Page: 3 of 28

Risk Characterization—Risk characterization combines chemical-specific toxicity information with quantitative and qualitative information from the exposure

assessment. This information is compared to measured contaminant exposure levels predicted through environmental fate and transport modeling. These comparisons

are used to determine whether concentrations of contaminants at or near the sites are

affecting, or could potentially affect, human health. An uncertainty analysis is also included in this section which presents critical assumptions, such as background

concentrations and conditions, that are considered uncertainties in this report.

9.2 CONTAMINANT IDENTIFICATION

In this section, the available information on the hazardous substances at the site is evaluated, the

chemicals detected in the environmental media (i.e., groundwater, sediment, and surface water)

sampled at the Evergreen Manor site are summarized, and the COPCs are identified.

9.2.1 Contaminant Characterization

Modio investigated during the Evergreen Manor Remedial Investigation (RI) included groundwater

sediment, and surface water. The following is a summary of the investigation results that are

described in more detail in Section 5. Figures 4-1 and 4-2 present the groundwater sample

locations. Figure 4-3 presents the sediment and surface water sample locations.

The geology underlying the site was characterized using cone penetration testing (CPT) methods.

Groundwater quality and flow conditions were assessed to evaluate areal and vertical extent of

contaminant migration and to determine the concentrations of the contaminants in groundwater.

Depth to water in monitoring wells was measured to determine the direction of groundwater flow

and hydraulic gradient. Residential wells, monitoring wells, and CPT boreholes were sampled to

determine groundwater quality within the unconfined sand and gravel aquifer.

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RFW036-2A-AHVH

Section: 9 Revision: 1

Date: 28 March 2001

Page: 4 of 28

Twenty-two groundwater samples were obtained from residential wells to determine the extent of

contamination and the concentrations of contaminants in the aquifer. All residential well samples

were analyzed for VOCs. In addition, six of the residential wells were analyzed for orthophosphate.

ammonia, nitrate, nitrite, COD, sulfide, and sulfate parameters to evaluate natural attenuation

characteristics across the lateral extent of the expected plume. Groundwater samples were collected

at 10 CPT locations to determine if groundwater has been impacted by contamination and to evaluate

the lateral and vertical extent of contamination. Discreet samples were taken at approximately 10

foot intervals for the purpose of determining the zone containing the maximum concentrations of

VOC contamination. Fifteen monitoring wells previously installed by IEPA were also sampled for

VOCs. Eight of these 15 monitoring wells were sampled for natural attenuation parameters to

determine groundwater quality and aid in determining the extent of contamination.

Six sediment samples were collected during the field investigation -- three sediment samples were

collected from the Rock River and three sediment samples were collected from Dry Creek. The

sediment samples were collected to determine if site contaminants have been discharged or have

migrated into the river and creek. The sediment samples were collected in areas that sediments are

readily deposited and were collected within approximately 5 feet of the shore. The six investigative

samples from Rock River and Dry Creek were submitted for VOC and TOC analysis.

Six surface water samples were collected during the field investigation -- three surface water samples

were collected from the Rock River and three surface water samples were collected from Dry Creek.

The surface water samples were collected to determine if site contaminants were migrating off-site

via surface water runoff or being discharged from groundwater. Surface water samples were

collected prior to collection of sediment samples at approximately the same locations as sediment

samples. Surface water samples were collected within 5 feet of the shore and just below the water

surface. The sampling progressed from downstream to upstream locations to minimize the impact

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Section: 9 Revision: 1

Date: 28 March 2001

Page: 5 of 28

of sediment disturbance and/or cross contamination of samples. The six investigative samples were

submitted for VOC analysis.

9.2.2 Data Evaluation

Chemical analyses were performed in a mobile field laboratory operated by the ESAT Region V

Mobile Laboratory. A stationary laboratory operated by Chemtech of Englewood, New Jersey, a

Contract Laboratory Program (CLP) laboratory, performed confirmatory CLP laboratory analysis

on 10 percent of the water and sediment VOC samples. All analyses were performed according to

the U.S. EPA-approved Quality Assurance Project Plan (QAPP) developed for the Evergreen Manor

site (WESTON, 2000a). All CLP generated data was validated by the U.S. EPA Region V Superfund

Division Field Services Section Quality Assurance Reviewer. In addition, WESTON's data

reviewers conducted a systematic review of the data for compliance with established QC criteria.

All SAS analytical data was also reviewed and validated by WESTON. The reader is referred to the

RI/FS Work Plan and QAPP (WESTON, 2000b and 2000a) for detailed information on data quality.

All environmental samples (i.e., groundwater, surface water, and sediment) collected during the Site

Characterization were analyzed for VOCs. Field duplicate samples were collected at selected

locations at a one per 10 frequency. The chemicals found in each environmental medium are

summarized by frequency of detection (i.e., the ratio of the number of samples in which the chemical

was detected to the number of samples available) and the minimum and maximum detected

concentrations in Table 2, "Occurrence, Distribution, and Selection of Chemicals of Potential

Concern," provided in Appendix A.

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Section: 9 Revision: 1

Date: 28 March 2001

Page: 6 of 28

9.2.3 Identification of Constituents of Potential Concern

The quantitative assessment of exposure, and consequently risk, for a site is based on those

chemicals considered as COPCs for the site. The COPCs are a subset of all the chemicals positively

identified at a site and are those constituents associated with site processes and measured above

background levels. The risks associated with the COPCs are expected to be more significant than

those associated with the other less toxic and less prevalent chemicals at the site. The list of COPCs

evaluated in a human health risk assessment may not be the same as that evaluated in an ecological

risk assessment.

Chemical COPCs identified at the Evergreen Manor site are VOCs. In general, the list of chemical

COPCs evaluated in the human health risk assessment includes those that are:

Positively detected in at least one CLP sample in a given medium, including: (a) chemicals with no qualifiers attached (excluding samples with unusually high

detection limits), and (b) chemicals with qualifiers attached that indicate known

rachitiles out alixilown or estillated concentrations (e.g., J-quarities data).

Detected at levels significantly elevated above levels of the same chemicals detected

in associated blank samples.

Detected at levels above Region 9 soil risk-based screening levels (U.S. EPA,

1999b), which are associated with a cancer risk of 1E-06 (one-in-one-million) and

a systemic hazard quotient (HQ) of 1. In order to provide a more conservative screening and to account for similar toxic endpoints among noncarcinogenic

compounds, a HQ of 0.1 was used in screening noncarcinogenic chemicals and a risk

level of 1E-07 was used in screening carcinogenic chemicals, based on U.S. EPA

(U.S. EPA 1993b) guidance. Where risk-based concentrations are available for cancer and non-cancer endpoints and both ingestion and inhalation exposure routes,

the lower (i.e., most stringent) value was used for the screening comparison.

The selection of COPCs for each environmental medium evaluated at Evergreen Manor is presented

in Table 2, "Occurrence, Distribution, and Selection of Chemicals of Potential Concern," which is

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Section: 9 Revision: 1

Date: 28 March 2001 Page: 7 of 28

provided in Appendix A. Since VOCs were not detected above method detection limits in surface

water at locations adjacent and downgradient of the site, surface water is not evaluated further in this

risk assessment.

VOCs were detected in sediment at several locations adjacent to the site; however, the constituent

detected, methyl acetate, does not have any established toxicity data. The only other VOCs detected

in sediment were detected at locations upgradient of the site. Therefore, the sediment medium is not

evaluated further in the risk assessment.

9.3 EXPOSURE ASSESSMENT

The purpose of the exposure assessment is to estimate the magnitude of human exposure to the

chemicals found in environmental media at the Evergreen Manor site. The results of the exposure

assessment are subsequently combined with the chemical-specific toxicity information to

quantitatively estimate the human health risks associated with chemical exposure at this site. The

identification of actual or potential pathways through which human receptors could be exposed to

chemicals in groundwater at the site includes identification and characterization of the site and the

potentially exposed populations. Exposure to surface water and sediment in the Rock River and Dry

Creek were not evaluated further in this risk assessment because contaminants were not detected in

surface water, contaminants were detected at upgradient sediment sampling locations only, or

toxicity data is not available for the contaminant detected in sediment adjacent to the site. After

exposure pathways have been identified, daily intakes of the COPCs are quantified using standard

exposure algorithms.

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Section: 9 Revision: 1

Date: 28 March 2001 Page: 8 of 28

9.3.1 Characterization of Exposure Setting

The first step in evaluating exposure is to characterize the site with respect to its physical

characteristics as well as the human populations on and near the site. Information gathered during

this step will support the identification of exposure pathways and the determination of exposure

assumptions.

The area in the vicinity of the site consists of a mixture of land uses including residential.

commercial, and light and heavy industrial. Based on the 1990 census, there are 2,079 people living

within the Village of Roscoe. The site includes four residential subdivisions and is located

approximately 1.5 miles northwest of the Roscoe city limits. Nearby development includes Roscoe

Rock & Sand, Inc., a gravel pit and concrete mixing facility, to the northeast of the site, and an

industrial park approximately 2 miles to the northeast of the site. The Hononegah Forest Preserve

is located to the west and agricultural fields to the north and east of the site. Additional information

on the physical setting of the site and surrounding area including climate, vegetation, soil type,

surface hydrology, and groundwater hydrology is presented in Section 3.

The Evergreen Manor site includes four residential subdivisions. Based on current site conditions

and site ownership, the baseline risk assessment evaluates residents as the current receptor group at

this site. The risk assessment also evaluated future residential and commercial/industrial use of the

site. Thus, potential risks were evaluated for current/future residential and future

commercial/industrial receptors.

9.3.2 Identification of Exposure Pathways

An exposure pathway generally consists of four elements: (1) a source and mechanism of

contaminant release, (2) a retention or transport medium, (3) a point of potential human contact with

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Section: 9 Revision: 1

Date: 28 March 2001

Page: 9 of 28

the contaminated medium (referred to as the exposure point), and (4) an exposure route (i.e.,

ingestion, dermal contact, or inhalation) at the exposure point (U.S. EPA, 1989). Table 1, "Selection

of Exposure Pathways," provided in Appendix A integrates and summarizes the information

concerning source areas, chemical migration pathways, receptor populations, and exposure routes

into a combination of potential human exposure pathways for the Evergreen Manor site. The

following subsections describe the process used to identify and select exposure pathways for

quantitative analysis.

Source Area

As previously discussed, the Evergreen Manor site includes four residential subdivisions. The site

area was used as farmland prior to development. The Hononegah Heights subdivision was developed

between 1940 and 1964; the Tresemer subdivision was developed between 1972 and 1974; the Olde

Farm subdivision was developed between 1976 and 1979; and the Evergreen Manor subdivision was

developed between 1986 and 1988. The source area was identified in an Action Memorandum (U.S.

EPA, 1999a) as a small industrial/commercial area at the intersection of Rockton Road and Illinois

Route 251. The specific source(s) of the contamination has not been discovered. The IEPA has

identified four potentially responsible parties based on sampling results and historic operations.

**Chemical Migration Pathways** 

This subsection briefly addresses the fate of chemicals measured in groundwater at the site, their

potential to be transported to other environmental media, and their potential to migrate off site. After

a chemical is released into the environment, it may be transported (e.g., advected downstream in

water), physically transformed (e.g., volatilized), chemically transformed (e.g., oxidation/reduction),

biologically transformed (e.g., biodegradation), or bioaccumulated in one or more media (U.S. EPA,

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Section: 9 Revision: 1

Date: 28 March 2001 Page: 10 of 28

1989). A goal of fate analysis is to identify other (non-source) environmental media and off-site

areas potentially affected by chemical migration.

Several VOCs were measured above screening levels in groundwater. These VOCs could migrate

toward downgradient receptor areas or into other environmental media (e.g., the Rock River or Dry

Creek). No VOCs were measured in surface water. VOCs measured in sediment were either at

upgradient locations, or did not have toxicity data. As the VOCs in groundwater reach a surface

waterbody, they are expected to readily volatilize to the atmosphere. However, volatile

contaminants in groundwater that are used as a household water supply can readily enter the

enclosed atmosphere of a residence during dishwashing, clothes laundering, and showering.

**Exposure Points and Exposure Routes** 

Two receptor groups were assumed to be exposed to constituents in environmental media at the site:

• Current/future on-site resident (adult and young child)

• Future commercial/industrial worker (adult)

The site is currently occupied by residences and surrounded by various commercial and industrial

businesses. Residents (adults and young children) and commercial/industrial workers from nearby

businesses use groundwater as their potable water supply. These receptor groups may potentially be

exposed to COPCs in groundwater. There are three primary exposure routes for chemicals in water:

ingestion, dermal absorption, and inhalation. Inhalation of volatile chemicals is considered routinely

only for chemicals with a Henry's Law constant of 1 x 10<sup>-5</sup> atm-m<sup>3</sup>/mole or greater and a molecular

weight of less than 200 g/mole (U.S. EPA, 1991a).

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Section: 9 Revision: 1

Date: 28 March 2001

Page: 11 of 28

**Exposure Pathways** 

For each receptor group, the following exposure pathways were quantitatively evaluated in the risk

assessment:

Ingestion of contaminated groundwater.

• Dermal absorption of chemicals from groundwater.

Inhalation of volatiles from groundwater.

9.3.3 Quantification of Exposure

The degree of receptor exposure that occurs through each exposure pathway is determined by

behavioral, chemical, and physiological factors. Behavioral factors affecting exposure would include

the amount of time spent on-site, the activities engaged in while on-site, and the amount and type

of clothing worn. Chemical factors affecting the degree of exposure include the extent to which a

chemical is absorbed through the skin and gastrointestinal tract (i.e., the absorption efficiency).

Physiological factors affecting exposure would include the ability of the body to metabolize and

eliminate the chemical(s). To quantify exposures in the risk assessment process, it is necessary to

make assumptions concerning these factors in the absence of specific, detailed information. These

assumptions are represented by a series of exposure parameters that quantify the magnitude,

frequency, and duration of the exposure. In addition, the quantification of exposure requires

estimates of chemical concentrations to which the receptor is exposed.

**Exposure Point Concentrations** 

The exposure point concentration is the concentration of a chemical to which a receptor may be

exposed. The exposure point concentration for each chemical in each medium is intended to

represent a reasonable maximum estimate of the concentration a receptor is likely to be exposed to

over time. Groundwater is generally evaluated at the center of any recognizable plume(s) of COPCs

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Section: 9 Revision: 1

Date: 28 March 2001

Page: 12 of 28

in potential sources of groundwater. A distinct plume was not recognizable at this site; therefore,

the maximum concentration of each COPC was used as the exposure point concentration for

groundwater.

Exposure point concentrations for groundwater water are summarized in Table 3, "Medium-Specific

Exposure Point Concentration Summary", which is provided in Appendix A.

Reasonable Maximum and Central Tendency Exposure

To evaluate exposures over the range of possible conditions that may exist at the Evergreen Manor

site, two hypothetical degrees of exposure are considered in this study following U.S. EPA (U.S.

EPA, 1992c; 1998) guidance. These degrees of exposure are reasonable maximum exposure (RME)

and central tendency (CT). The RME is the highest exposure that is reasonably expected to occur

at a site; the CT is intended to represent more typical (i.e., central tendency) exposure conditions.

In evaluating RME and CT scenarios, the exposure point concentration remains the same, while the

exposure parameters are adjusted to reasonable maximum and central tendency values.

**Exposure Algorithms** 

U.S. EPA has developed exposure algorithms for use in calculating chemical intakes through the

exposure pathways and routes that are relevant for this site. These algorithms combine chemical

exposure point concentrations with pathway- and route-specific parameters to produce daily

chemical intakes in terms of the milligrams of chemical taken into the body per kilogram of body

weight per day (mg/kg-day). The following subsections discuss the exposure pathways and routes

through which receptors are assumed to be exposed to site contaminants, and present the exposure

algorithms and exposure parameters that were used in this risk assessment.

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Section: 9 Revision: 1

Date: 28 March 2001 Page: 13 of 28

The basic mathematical models used to calculate intakes are presented in the following subsections.

Each model defines the exposure variables used in estimating the intake, and includes the

assumptions (e.g., exposure parameters) used in the model. In general, the exposure parameters used

are standard values recommended by U.S. EPA guidance (U.S. EPA, 1989; 1991c; 1997). When

available, site-specific exposure information was used.

Water Pathway

The current/future resident and future commercial/industrial worker may be exposed to groundwater

used as a potable water supply The estimated daily intakes that result from groundwater exposure

are presented in Appendix A.

Ingestion of Water

Chemical intake through water ingestion and dermal absorption is calculated as follows:

Intake =  $C_w \times IR_w \times EF \times ED/(BW \times AT)$ 

where:

Intake = Estimated water intake (mg/kg-day)

 $C_w = Chemical concentration in water (mg/L)$ 

ED = Exposure duration (yrs)

EF = Exposure frequency (days/yr)

 $IR_w = Ingestion rate (L/day)$ 

BW = Body weight (kg)

AT = Averaging time (days)

Dermal Absorption from Water

Intake =  $C_w x SA x PC x 1L/1000 cm^3 x ET x EF x ED/ (BW x AT)$ 

where:

Intake = Estimated water intake (mg/kg-day)

 $C_w$  = Chemical concentration in water (mg/L)

SA = Surface area available for contact (cm<sup>2</sup>)

Section: 9 Revision: 1

Date: 28 March 2001 Page: 14 of 28

```
PC
              =
                     Permeability constant (cm/hr)
                     Exposure time (hours/day)
ET
              =
                     Exposure frequency (days/yr)
EF
                     Exposure duration (yrs)
ED
              =
                     Ingestion rate (L/day)
IRw
              =
                     Body weight (kg)
BW
              =
                     Averaging time (days)
AT
              =
```

Inhalation of Organic Vapors from Groundwater

For organic compounds detected in groundwater, vapor inhalation intake is calculated as:

```
Intake =
                        C_w \times EF \times IH \times K \times ED / (BW \times AT)
where:
       Intake
                                Estimated daily intake via vapor inhalation (mg/kg-day)
                               Chemical concentration in water (mg/L)
        C<sub>w</sub>
                        =
        ED
                        =
                               Exposure duration (yrs)
                               Exposure frequency (days/yr)
       EF
                       =
                               Inhalation rate (m³/day)
       IH
                       =
                               Volatilization factor (unitless)
       K
       BW
                               Body weight (kg)
                       =
        AT
                               Averaging time (days)
```

### **Exposure Parameters**

Values used for daily intake calculations are presented in Table 4, "Values Used for Daily Intake Calculations," and are provided in Appendix A.

Residential and commercial/industrial worker exposure to groundwater through drinking water ingestion, dermal absorption, and inhalation were considered. Exposures through drinking water may occur by a variety of mechanisms, including ingestion of drinking water and ingestion of foods prepared with or in water. The reasonable maximum ingestion rate for potable water is established at 2 L/day for an adult and 1 L/day for a child (U.S. EPA, 1989). The average drinking water ingestion rate is 1.4 L/day for an adult and 0.7 L/day for a child (U.S. EPA, 1997). For commercial/industrial workers, the reasonable maximum water ingestion rate is 1 L/day while the

Section: 9 Revisión: 1

Date: 28 March 2001

Page: 15 of 28

average ingestion rate is 0.7 L/day. These amounts include water consumed in the form of other

beverages and the ingestion of foods prepared in or with water.

Certain non-potable uses of water may result in skin contact and dermal absorption of waterborne

contaminants. For bathing and swimming scenarios, 75 to 100% of the skin surface is exposed. The

total adult body surface area can vary from 17,000 to 23,000 cm<sup>2</sup> for an adult. A mean value of

20,000 cm<sup>2</sup> and a reasonable maximum value of 23,000 cm<sup>2</sup> are recommended as default values by

U.S. EPA (1992a). The total body surface area for a 1 to 6-year old child is 8,538 cm<sup>2</sup> for the 95th

percentile and 7,314 cm<sup>2</sup> for the 50<sup>th</sup> percentile (U.S. EPA, 1997). Shower times have been found

to range from 10 to 35 minutes (U.S. EPA, 1997). The permeability constants for the organic

COPCs are listed in Table 9-1.

Indoor inhalation of volatiles is relevant only for chemicals that easily volatilize, i.e., chemicals with

a Henry's law constant greater than 10-5 atm-m<sup>3</sup>/mole and a molecular weight of less than 200

g/mole. Andelman (1990, as cited in U.S. EPA, 1991c), derived an equation that relates the

concentration of a contaminant in household water and the average concentration of the volatilized

contaminant in air. The volatilization constant is 0.0005 x 1000 L/m<sup>3</sup>, where the 1000 L/m<sup>3</sup>

conversion factor is used so that the resulting air concentration is expressed as mg/m<sup>3</sup>. The indoor

air inhalation rate for an adult is 15 m<sup>3</sup>/day (U.S. EPA, 1991c) and the indoor inhalation rate for a

young child is 8.7 m<sup>3</sup>/day (U.S. EPA, 1997).

An exposure frequency (EF) of 350 days per year is the default reasonable maximum value for

residents; it was also used for the average scenario. This EF is based on the common assumption that

workers take two weeks of vacation per year to support a value of 15 days per year spent away from

home (U.S. EPA, 1991a). An EF of 250 days per year is the default value for commercial/industrial

workers for both the RME and CT scenarios. Again, it is assumed that a worker takes two weeks of

vacation per year (U.S. EPA, 1991a).

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Section: 9 Revisión: 1

Date: 28 March 2001 Page: 16 of 28

The national upperbound time at one residence is 30 years (U.S. EPA, 1991c) and the average time

at one residence is 9 years (U.S. EPA, 1997). The exposure duration for an adult was adjusted to

24 years for the RME scenario and 7 years for the CT scenario to account for differences in ingestion

rate and body weight between children and adults. These values were applied for the adult resident.

The reasonable maximum ED for a 1- through 6-year-old child is 6 years; an average exposure

duration of 2 years was used, which was estimated based on the ratio of the average to maximum

adult exposure duration. For the commercial/industrial worker, the ED is assumed to be an average

of 9 years (U.S. EPA, 1992b) and a maximum of 25 years (U.S. EPA, 1991c).

The value for body weight (BW) is the average body weight over the exposure period. An average

BW is used because, when combined with other variable values, it is believed to result in the

reasonable maximum exposure. Incorporating a higher BW with the same intake rate would result

in lower exposure than the reasonable maximum. In addition, using an average BW rather than a

reasonable maximum is recommended because the available toxicity data are based on average body

weight. The recommended average BW for an 18- to 75-year old adult is 70 kg (U.S. EPA, 1991c).

The recommended average BW for a 1- through 6-year-old is 15 kg (U.S. EPA, 1991c). An average

BW is applied for both the RME and CT scenarios.

The averaging time (AT) is the period over which exposure is averaged. For non-carcinogenic

effects, AT is equal to exposure duration (ED). For carcinogens, AT is equal to a 70-year lifetime

(U.S. EPA, 1991c).

9.4 TOXICITY ASSESSMENT

A toxicity assessment presents the appropriate toxicity values and the weight of evidence for the

toxicity of each of the COPCs. Applicable human toxicity values are identified for each COPC for

the relevant exposure routes. These toxicity values include reference doses (RfDs) for evaluating

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Section: 9 Revision: 1

Date: 28 March 2001 Page: 17 of 28

potential non-carcinogenic health effects and cancer slope factors (CSFs) for evaluating carcinogenic

risks.

Toxicity criteria were obtained from the following sources:

• Integrated Risk Information System (IRIS) (U.S. EPA, 2000)

Health Effects Assessment Summary Tables (HEAST) (U.S. EPA, 1995)

If toxicity criteria were not available from these sources, toxicity criteria presented in the Region 9

Risk-Based Concentration Tables (U.S. EPA, 1999b) were used. The toxicity criteria for the COPCs

are presented in Appendix A (Tables 5-1 and 5-2 for noncarcinogenic effects and Tables 6-1 and

6-2 for carcinogenic effects). Chemicals of potential concern lacking toxicity criteria will be

discussed in the Uncertainty Analysis.

9.5 RISK CHARACTERIZATION

In a risk characterization, the results of the exposure assessment and the toxicity assessment are

integrated to quantitatively evaluate the potential current and future risk to human health.

Carcinogenic and noncarcinogenic risks are evaluated for each COPC through each exposure route

of concern and for all COPCs through all exposure routes combined. The risk characterization also

identifies uncertainties associated with contaminant, toxicity, or exposure assumptions. Tables

containing the risk calculations following the Risk Assessment Guidance for Superfund, Part D (U.S.

EPA, 1998) format are provided in Appendix A.

9.5.1 Quantitative Evaluation for Chemical COPCs

Non-carcinogenic effects are evaluated by comparing estimated daily intakes of chemical COPCs

to RfDs. This is accomplished by calculating hazard quotients (HQs) and hazard indices (HIs). An

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Revisión: I

Date: 28 March 2001 Page: 18 of 28

HQ for a particular COPC through a given exposure route is the ratio between the estimated daily intake and the applicable RfD, as shown in the following equation:

HQ = EDI / RfD

where:

HQ = Hazard quotient (unitless)

EDI = Estimated daily intake (mg/kg-day)

RfD = Reference dose (mg/kg-day)

Screening level HIs were calculated by summing HQs across all exposure pathways and all target effects for all COPCs. Separate HIs were calculated for child and adult receptors. More appropriately, an HI is developed for chemicals that affect the same target organ or produce the same critical effect since the assumption of dose additivity is most properly applied to compounds that induce the same effect by the same mechanism of action (U.S. EPA, 1989). This approach is based on the assumption that combined exposure to several chemicals below their threshold level can result in an adverse health effect when they have the same critical effect or the same target organ. If the screening level HI exceeded one, chemical COPCs were segregated by target organ and a separate HI value for each effect/target organ was calculated. If the HQ for any chemical or if the HI value for any effect/target organ exceeds one, non-carcinogenic health effects are considered possible.

Carcinogenic risks were calculated for each carcinogen through each exposure pathway for each receptor. In risk assessment calculations, cancer risks are estimated as the incremental, or excess probability of an individual developing cancer over a lifetime as a result of exposure to a carcinogen. This risk is in addition to the lifetime cancer risk experienced by the general, non-exposed population. Cancer risks were calculated for each chemical COPC using the following formula:

Risk = EDI x CSF

where:

Risk = Excess cancer risk (unitless probability)

EDI = Estimated daily intake (mg/kg-d) (see Section 3.3.2)

Section: 9 Revision: 1

Date: 28 March 2001 Page: 19 of 28

CSF = Cancer slope factor  $(mg/kg-d)^{-1}$ 

The total risk posed by each chemical COPC will be calculated by adding risks posed by the COPC

through all exposure routes. The lifetime incremental cancer risk posed by all chemical COPCs is

estimated by summing the adult and child risks posed by all chemical COPCs through all exposure

routes.

9.5.2 Residential Scenario

Under this current/future exposure scenario, it was assumed that residents use on-site groundwater

as a potable water supply, potentially being exposed to chemical COPCs in groundwater through

ingestion, dermal absorption while bathing, and inhalation of volatiles. Details of all risk

calculations for this receptor group are presented in Appendix A.

Cancer Risk

The potential chemical cancer risk estimates associated with the residential scenario are presented

in Table 9-2. The chemical cancer risk ranged from 4.6E-06 to 1.9E-05. Tetrachloroethene had an

individual RME cancer risk estimate exceeding 1.0E-06 via ingestion, while trichloroethene,

tetrachloroethene, and chloroform had individual RME cancer risk estimates greater than 1.0E-06

via inhalation.

Non-Cancer Health Effects

The estimates of the potential for adverse non-carcinogenic health effects associated with the

residential scenario are presented in Table 9-3. The total HI (all COPCs, all target effects, all

exposure routes) ranged from 1.4 to 1.4 for the adult resident and from 3.7 to 3.8 for the child

resident. Chloroform was the only COPC with an individual HQ value (via inhalation) exceeding

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RFW036-2A-AHVH

Section: 9 Revision: 1

Date: 28 March 2001 Page: 20 of 28

one. Acetone, tetrachloroethene, and methylene chloride effect the same target organ (liver) as

chloroform. These COPCs with the same target organ/effect had a total HI (based on effect) greater

than one only when combined with chloroform.

9.5.3 Commercial / Industrial Worker Scenario

Under this exposure scenario, it is assumed that businesses are developed within the Evergreen

Manor site. Commercial/industrial receptors were assumed to use on-site groundwater as a potable

water supply, potentially being exposed to chemical COPCs in groundwater through ingestion,

dermal absorption while bathing, and inhalation of volatiles. Details of all risk calculations for this

receptor group are presented in Appendix A.

Cancer Risk

For this future groundwater pathway, the chemical cancer risk ranged from 2.0E-06 to 6.9E-06.

Tetrachloroethene had an individual RME cancer risk estimate exceeding 1.0E-06 via ingestion and

dermal absorption, while chloroform had an individual RME cancer risk estimate exceeding 1.0E-06

for inhalation.

**Non-Cancer Health Effects** 

For the future groundwater pathway, the total HI ranged from 0.97 to 0.99. No COPCs with the

same target organ/effect had a total HI (based on effect) greater than one in any of groundwater

samples.

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Section: 9 Revision: 1

Date: 28 March 2001 Page: 21 of 28

9.5.4 Uncertainty Analysis

The goal of an uncertainty analysis in a risk assessment is to provide to the appropriate decision

makers (i.e., risk managers) a wide range of information about the key assumptions, their inherent

uncertainty and variability, and the impact of this uncertainty and variability on the estimates of risk.

The uncertainty analysis should show that risks are relative in nature and do not represent an

absolute quantification. This is an important point that is vital to the proper interpretation and

understanding of the risks presented in this report. This subsection attempts to explain the key

assumptions used in this risk assessment and present a range of risks covering the variability inherent

in these assumptions.

There are three areas in this report with significant levels of uncertainty, which are described in the

following subsections:

• Environmental data used in risk assessment.

Exposure assumptions.

Toxicological assumptions.

**Environmental Data Used in Risk Assessment** 

The environmental data collected from a site and how it is used in a risk assessment contributes

uncertainty to the risk estimates. There is a measure of uncertainty associated with the exposure

point concentrations used for the groundwater COPCs. The exposure point concentrations were set

at the maximum detected concentration within the plume, and does not consider the possibility that

some parts of the site might be more or less contaminated than others, and therefore the actual

exposure point concentration might be different than the calculated values.

**Exposure Assumptions** 

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Section: 9 Revision: 1

Date: 28 March 2001

Page: 22 of 28

There are a number of exposure-related assumptions that are likely to result in significant

uncertainty. In most cases, this uncertainty overestimates the realistic exposures, and therefore,

overestimates the risk. This is appropriate when performing risk assessments of this type so that the

risk managers can be reasonably assured that the risks to the public are not underestimated, and so

that risk assessments for different locations and different scenarios can be compared.

Much of the uncertainty involves the use of standard exposure factors relating to a drinking water

ingestion rate, frequency of exposure, etc. These factors are designed to cover reasonably maximum

exposed individuals who are at the site for many years. It is very likely that an actual individual

would be exposed to a lesser degree than the reasonable maximum, and possibly to a significantly

lesser degree. Specific exposure assumptions that add uncertainty to the risk estimates are described

below.

**Toxicity Assessment** 

Toxicological uncertainties primarily relate to the methodologies by which carcinogenic and non-

carcinogenic health criteria are developed. Standard slope factors and reference doses established

by the U.S. EPA were used to estimate potential carcinogenic and non-carcinogenic health effects

from exposure to chemical COPCs at the site. Cancer slope factors are derived using a non-threshold

theory of cancer formation, which assumes there is no "safe" level of exposure to any carcinogen.

The animal studies from which slope factors are usually derived are performed at high doses. The

dose-response data from these studies is typically extrapolated down many orders of magnitude to

estimate risks associated with the comparatively low dose to which humans might be exposed

through environmental contamination. This low dose extrapolation produces conservative estimates

of risk, possibly to a very significant degree. However, based on the current state of knowledge, this

level of conservatism cannot be quantified.

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Section: 9 Revision: 1

Date: 28 March 2001

Page: 23 of 28

There is also uncertainty associated with the estimation of non-cancer risks. In this case, there is an

assumed threshold that can be tolerated by any individual without the threat of a potential health

impact. The approach is to identify a sub-threshold level (i.e., the reference dose) that will be

protective of the most sensitive individuals in the population. The reference dose is usually based

on animal studies or limited human data and incorporates uncertainty factors anywhere from one to

five orders of magnitude. These factors reflect the degree of extrapolation used in the derivation of

a reference dose (U.S. EPA, 1989).

The application of a route-specific slope factor and reference dose to another exposure route

contributes uncertainty to the evaluation of risks. Most reference doses and some slope factors are

expressed as the amount of a substance administered, while dermal exposure estimates are expressed

as an absorbed dose. Ideally, each oral toxicity value should be adjusted by an appropriate

gastrointestinal absorption factor (U.S. EPA, 1989). The lower the gastrointestinal absorption factor,

the more conservative the derived toxicity values. There is limited information on the absorption

of the COPCs: thus, no adjustments were made to the oral toxicity values. Assuming 100%

gastrointestinal absorption may lead to a non-conservative estimate of a dermal toxicity value.

However, the use of adequately conservative uncertainty factors in the derivation of toxicity values

is expected to ensure that the health criteria used are adequately protective of human health.

Summary

Table 9-4 presents a qualitative evaluation of the effects of each of these three key areas of

uncertainty on the estimation of risk for the Evergreen Manor site. The risks presented in this report

need to be viewed in light of the inherent uncertainty, which is summarized in this table. Column

1 lists the uncertainty elements identified as key contributors to this risk assessment. Column 2

shows the assumptions that represent a likely moderate to high overestimation of risk, while

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Section: 9 Revision: 1

Date: 28 March 2001 Page: 24 of 28

potential underestimations of risk are noted in Column 3. Column 4 identified several areas where

the potential exists for low to moderate over- or underestimation of risk.

In summary, the level of uncertainty in this risk assessment is moderate. Most of the uncertainty

results in overestimating risk, while some may result in either an over- or an underestimation of the

risk, and some may result in an underestimation of risk. However, it is likely that the overall risk

is overestimated by an order of magnitude.

Evergreen Manor Remedial Investigation Report Section: 9 Revision: 1

Date: 28 March 2001 Page: 25 of 28

Table 9-1

# Permeability Coefficients for Organic COPCs Evergreen Manor Site Roscoe, Winnebago County, Illinois

Chemical of Potential Concern	Permeability Coefficient (cm/hr)
Acetone	5.7E-04
Benzene	2.1E-02
Methylene chloride	4.5E-03
Tetrachloroethene	4.8E-02
Trichloroethene	1.6E-02
Chloroform	8.9E-03

Section: 9 Revision: 1

Date: 28 March 2001 Page: 26 of 28

Table 9-2

# Total Carcinogenic Risk Associated with Chemical COPC Exposure Evergreen Manor Site Roscoe, Winnebago County, Illinois

	Total Lifetime Cancer Risk							
Exposure Route		l Land Use +Adult)	Commercial/Industrial Land Use (adult)					
	RME	CT	RME	CT				
Ingestion	8.6E-06	1.8E-06	2.0E-06	5.1E-07				
Dermal absorption	2.2E-06	1.7E-07	1.1E-06	1.0E-07				
Inhalation	8.5E-06	2.6E-06	3.8E-06	1.4E-06				
Subtotal	1.9E-05	4.6E-06	6.9E-06	2.0E-06				

Evergreen Manor Remedial Investigation Report Section: 9 Revision: 1

Date: 28 March 2001 Page: 27 of 28

Table 9-3

# Total Hazard Index Associated with Chemical COPC Exposure Evergreen Manor Site Roscoe, Winnebago County, Illinois

	Total Hazard Index									
Exposure Route	Residential Land Use (Child)		Residential Land Use (Adult)		Commercial/Industrial Land Use (Adult)					
<u> </u>	RME	CT	RME	CT	RME	CT				
Ingestion	2.0E-01	1.4E-01	8.8E-02	6.1E-02	3.1E-02	2.2E-02				
Dermal absorption	2.1E-02	5.2E-03	1.2E-02	3.0E-03	8.5E-03	2.2E-03				
Inhalation	3.6	3.6	1.3	1.3	9.5E-01	9.5E-01				
Subtotal	3.8	3.7	1.4	1.4	9.9E-01	9.7E-01				

Section: 9 Revision: 1

Date: 28 March 2001 Page: 28 of 28

# Table 9-4

# Summary of Uncertainty Analysis Evergreen Manor Site Roscoe, Winnebago County, Illinois

	Effect on Risk Estimates				
Uncertainty Element	Potential for Overestimation	Potential for Underestimation	Potential for Over- or Underestimation		
Environmental Data					
Insufficient data to characterize media being evaluated	·		Low		
Systematic or random errors in the chemical analyses yielding erroneous data			Low		
Elimination of chemicals from quantitative analysis based on background levels and risk-based screening		Low			
Use of current exposure concentrations to represent future conditions (i.e., assumption of no attenuation of site chemicals)	Moderate		·		
Exposure Parameter Estimation					
Standard assumptions regarding body weights,			Low		
expectancy					
Media intake rates	Moderate				
Dermal absorption factors	Moderate				
Toxicity Data					
Use of U.S. EPA RfDs/SFs	Moderate-High				
Use of oral toxicity criteria for dermal exposure		Low			

Section: 10 Revision: 1

Date: 28 March 2001 Page: 1 of 13

### **SECTION 10**

## ECOLOGICAL RISK ASSESSMENT

A preliminary or screening-level risk evaluation is the initial ecological risk assessment screening at a hazardous waste site (U.S. EPA, 1996d). A screening level risk evaluation was conducted at the Evergreen Manor site to evaluate the potential impacts of contaminants on ecological receptors inhabiting the site and adjacent areas. U.S. EPA (1997b) defines a screening-level ecological risk assessment as "a preliminary risk assessment that can be conducted with limited site-specific data by defining assumptions for parameters that lack site-specific data." To ensure that sites which may pose an ecological risk are properly identified, U.S. EPA (1997b) suggests that "values should be consistently biased in the direction of overestimating risk. Without this bias, a screening evaluation could not provide a defensible conclusion for an absence of ecological risk." In conjunction with the human health risk assessment, the ecological risk assessment forms the basis for determining the need for remedial activities at a site and serves as the justification for the selected remedial action.

Technical risk assessment guidance for the performance of the screening-level ecological risk assessment came primarily from:

- Ecological Risk Assessment Guidance for Superfund: Process for Designing and Conducting Ecological Risk Assessments (U.S. EPA, 1997b).
- Guidelines for Ecological Risk Assessment (U.S. EPA, 1998b).

This screening-level ecological risk assessment consists of the following two steps:

- Screening-Level Problem Formulation and Screening-Level Ecological Effects Evaluation
- Screening-Level Preliminary Exposure Estimate and Risk Calculation

Section: 10 Revision: 1

Date: 28 March 2001 Page: 2 of 13

Each step of the screening-level ecological risk assessment for the Evergreen Manor site is presented

in the following sections.

10.1 SCREENING-LEVEL PROBLEM FORMULATION

The screening-level problem formulation step focuses on identifying categories of potential

ecological receptors that may exist in the site area; identifying contaminants which may pose

unacceptable risk to those receptors; and determining contaminant fate/transport and toxicity

mechanisms (U.S. EPA, 1996d). It is a planning step that identifies the major factors (i.e.,

environmental setting, extent of contamination, contaminant fate and transport, potential receptors,

and complete exposure pathways) to be considered in the screening-level ecological risk assessment.

The problem formulation is "the formal process of generating and evaluating preliminary hypotheses

about why ecological effects have occurred or may occur from human activities" (U.S. EPA, 1998b).

It is a planning step that identifies the major factors (e.g., site ecology, extent of contamination,

potential ecological receptors) to be considered in the assessment. The problem formulation

• Environmental setting and contamination.

Contaminant fate and transport.

Toxic mechanisms and potential receptors.

Complete exposure pathways.

General assessment endpoints.

Conceptual model.

10.1.1 Environmental Setting

10.1.1.1 Site Description

A description of the physical features of the Evergreen Manor site and the history of its use and past

removal activities are provided in Section 2 of this document. Further information on the physical

CHLAN01\WP\WO\RAC\036\29672-S10.WPD

RFW036-2A-AHVH

Evergreen Manor

Remedial Investigation Report

Section: 10 Revision: 1

Date: 28 March 2001

Page: 3 of 13

setting of the site and surrounding area including climate, vegetation, soil type, surface hydrology,

and groundwater hydrology is presented in the RI/FS Work Plan (WESTON, 2000b) and the Quality

Assurance Project Plan (WESTON, 2000a) and in Section 3 of this document.

10.1.1.2 Site Ecology

The site is located in the Central Lowland geomorphic province, in the eastern broadleaf forest

province of the Hot Continental Division in the Humid Temperate Domain (USDA Forest Service,

Ecological Subregions of the United States, http://www.fs.fed.us/land/pubs/ecoregions).

The Rock River receives drainage from three major streams - the Pecatonica River, the Kishwaukee

River, and the Green River. It is 163 miles long in Illinois, and drains 2,272,000 acres in Illinois.

Of the total river miles in this basin, 69 stream miles have "good" overall resource quality and 97.9

miles have "fair" quality. The Rock River enters the Mississippi River at Rock Island (IDNR,

http://dnr.state.il.us/lands/education/valerie/end/page6.htm). At Rockton, the mean daily discharge

ranges from 2820 afe in Santamber to 7375 afe in April with an annual mean of 1178 afe MISGS

CD-ROM, Current Year Discharge, http://www.il.water.usgs.gov/cd04-99/dis tbl/05437500.htm).

Dry Creek, a tributary of the Rock River, enters the river northwest of the Tresemer Subdivision.

West of Dry Creek, the river is classified by the NWI as a riverine wetland and east of the creek, the

river is classified as a lacustrine system. Forested wetlands border both the river and the creek west

of the site and the river south of the site. There are small areas of emergent wetlands within the

Evergreen Manor subdivision (Figure 10-1).

The U.S. Fish and Wildlife Service (USFWS) was contacted to obtain information on threatened and

endangered species within the Evergreen Manor project area. Species that may be present in the area

include the endangered Indiana bat (Myotis sodalis), the threatened prairie bush clover (Lespedeza

leptostachya), the threatened bald eagle (Haliaeetus leucocephalus). Further information on these

species and their critical habitat is provided in Appendix A.

CHLAN01\WP\WO\RAC\036\29672-S10.WPD

RFW036-2A-AHVH

Section: 10 Revision: 1

Date: 28 March 2001

Page: 4 of 13

10.1.2 Extent of Contamination

Information on the extent of contamination at the Evergreen Manor site is summarized in Section 7

of this document. Information on field sampling methods are presented in the Quality Assurance

Project Plan (WESTON, 2000a), the RI/FS Work Plan (WESTON, 2000b), and Section 4 of this

document. All data used in this evaluation were obtained during the RI. Elevated levels of VOCs

were measured in groundwater. VOCs were not detected in surface water, and downgradient of the

site, only one VOC was detected at a low concentration in sediment.

10.1.3 Contaminant Fate and Transport

While ecological receptors are generally not exposed to groundwater; the Rock River and Dry Creek

are adjacent to this site, and migration of groundwater contaminants to surface water is a potential

concern at this site; thus, this medium will be evaluated in the ecological risk assessment.

Information obtained during the site investigation indicated that contamination is restricted to the

unconfined and and aroust outwork aguifar. The water table is annovimately 35 feet has in this

aquifer.

10.1.4 Potential Ecological Receptors

Ecological receptors potentially exposed to groundwater contaminants from the Evergreen Manor

site include animals and plants common to rivers and streams of northwestern Illinois. A quantitative

survey of wildlife was not performed at this site as part of the screening-level ecological risk

assessment, though wildlife observations were made during RI activities. Aquatic biota potentially

inhabiting the Rock River and Dry Creek are the primary receptors at this site.

CHLAN01\WP\WO\RAC\036\29672-S10.WPD

RFW036-2A-AHVH

Section: 10 Revision: 1

Date: 28 March 2001

Page: 5 of 13

10.1.5 Complete Exposure Pathways

For an exposure pathway to be complete, a contaminant must be able to travel from the source to

ecological receptors and to be taken up by the receptors via one or more exposure routes (U.S. EPA,

1998b). For aquatic organisms, direct contact with water through the gills or dermis and ingestion

of water, food, and sediments are the primary exposure routes.

10.1.6 General Assessment Endpoints

Assessment endpoints are "explicit expressions of the environmental value that is to be protected"

(U.S. EPA, 1998b). The ecological resources selected to represent management goals for

environmental protection are reflected in the assessment endpoint. Assessment endpoints link the

risk assessment to management concerns and they are central to conceptual model development

(U.S. EPA, 1998b). The following principal criteria are used when selecting assessment endpoints

(U.S. EPA, 1998b):

The contaminants present and their concentrations.

• Mechanisms of toxicity of the contaminants to different groups of organisms.

Ecologically relevant receptor groups that are potentially sensitive or highly exposed

to the contaminant and attributes of their natural history.

Potentially complete exposure pathways.

The preliminary assessment endpoints for the Evergreen Manor site are changes in the aquatic

community structure and function attributable to COPCs measured in groundwater discharging to

surface water in the Rock River and Dry Creek.

CHLAN01\WP\WO\RAC\036\29672-S10.WPD

RFW036-2A-AHVH

Section: 10

Revision: 1 Date: 28 March 2001

Page: 6 of 13

10.1.7 Conceptual Model

The conceptual model establishes the complete exposure pathways that are evaluated in the

ecological risk assessment and the relationship of the measurement endpoints to the assessment

endpoints (U.S. EPA, 1997b). The conceptual model for the Evergreen Manor site is presented in

Table 10-1. Based on the conceptual site model, the exposure scenario included in the

environmental evaluation of the site was an aquatic community hazard quotient evaluation for fish

and other aquatic organisms that are directly exposed to COPCs in groundwater that discharges to

surface water, where media concentrations are compared with surface water quality benchmarks.

10.2 SCREENING-LEVEL ECOLOGICAL EFFECTS EVALUATION

In the ecological effects evaluation, information on the toxicity of the COPCs to ecological receptors

is presented. Toxicity information was used to develop toxicity reference values (TRVs) for selected

indicator species or communities. TRVs are expressed as an acceptable daily dose or media

annountration domanding on the recontaries

10.2.1 Constituents of Potential Concern

For groundwater, all VOCs detected above method detection limits are considered to be COPCs.

10.2.2 Toxicity Reference Values

For each COPC with a potentially complete exposure pathway, a screening-level toxicity reference

value (TRV) was developed from a review of literature. TRVs based on media concentrations are

not specific to individual species but instead are applicable to groups of organisms or communities

occupying the same medium (e.g., aquatic biota in surface water). The potential exists for ecological

receptors to indirectly contact groundwater through existing or potential discharge to sediments and

CHLAN01\WP\WO\RAC\036\29672-S10.WPD

RFW036-2A-AHVH

Remedial Investigation F Section: 10

Section: 10 Revision: 1

Date: 28 March 2001

Page: 7 of 13

surface water. Surface water TRVs are used to conservatively evaluate this potential exposure

pathway, assuming no attenuation or dilution of contaminants.

As a means of characterizing aquatic toxicity, national ambient water quality criteria (NAWQC)

have been developed for the protection of 95 percent of all aquatic life where sufficient data are

available (U.S. EPA, 1992e). Ambient water quality criteria are designed to be protective of all

aquatic biota occupying the same aquatic community or body of water. Not only fish, but also

aquatic invertebrates and plants are protected (U.S. EPA, 1986). However, NAWQC are not

available for the COPCs (U.S. EPA, 1999c). In addition, no state general use water quality standards

are available for the COPCs (35 IAC Part 302). Since water quality standards are not available for

the COPCs, alternative screening values are proposed. Tier II values described in the Proposed

Water Ouality Guidance for the Great Lakes System (U.S. EPA, 1993, as presented in Suter and

Tsao, 1996) and the lowest chronic value (LCVs) presented in the literature for a chemical (Suter

and Tsao, 1996) are proposed as surface water screening benchmarks.

10.2 CODERNING A DATE DADOCHDE ECHIMANTE

The screening-level exposure estimation involves the selection of exposure parameters for use in

calculating a daily exposure dose or exposure concentration. Measured environmental medium

concentrations (e.g., surface water, groundwater, sediment, and soil) are used for estimating exposure

of terrestrial and aquatic wildlife to site contaminants. Chemical exposure by aquatic life such as

fish are evaluated in this assessment through direct comparison to ecological benchmarks rather than

dose calculations. To estimate chemical exposure by aquatic life, groundwater concentrations are

compared directly to toxicity benchmarks (e.g., surface water screening benchmarks).

CHLAN01\WP\WO\RAC\036\29672-S10.WPD

RFW036-2A-AHVH

Section: 10 Revisión: 1

Date: 28 March 2001 Page: 8 of 13

10.4 <u>SCREENING-LEVEL RISK CHARACTERIZATION</u>

The screening-level risk characterization integrates information from the screening-level problem

formulation, screening-level ecological effects evaluation, and the screening-level exposure estimate

to predict the nature and extent of ecological risk or threat, as well as the environmental impact of

previous site activities.

A comparison of groundwater chemical concentrations with freshwater surface water screening

values was used to assess the likelihood of adverse effects to aquatic organisms inhabiting the Rock

River and Dry Creek, which are downgradient of the Evergreen Manor site (Table 10-2).

Comparison of groundwater chemical concentrations with surface water screening benchmarks were

used to evaluate potential for adverse impacts to aquatic organisms from contaminants discharging

to surface water. No screening benchmarks were exceeded; thus, negligible risk to aquatic receptors

is expected at this site.

10 11 Summar PTTmonutaints

The ecological risk assessment process is subject to a variety of uncertainties. Almost every step

involves assumptions based on professional judgment. Due to the conservative nature of a

screening-level ecological risk assessment, most of the uncertainty results are an overestimation of

risk. However, the risk may also be underestimated or unknown. Uncertainties specific to this risk

assessment can be attributed to:

• Environmental chemistry and sampling analysis.

Fate and transport parameters.

Exposure assumptions.

Toxicological data.

Uncertainties specific to this risk assessment include the following:

CHLAN01\WP\WO\RAC\036\29672-S10.WPD

RFW036-2A-AHVH

Section: 10 Revisión: 1

Date: 28 March 2001

Page: 9 of 13

 Maximum values were used as exposure point concentrations for all media. This is likely to result in an overestimation of risk, especially for aquatic ecological

receptors who may inhabit a greater portion of a stream than the area represented by

just one or a few samples

• Dilution and attenuation of VOCs in groundwater were not accounted for in

estimating future surface water concentrations in the river and creek. These fate and transport processes are expected to occur, resulting in lower discharge

concentrations.

• Tier II values were developed with fewer data than are required for the NAWQC.

The Tier II values are concentrations that would be expected to be higher than

NAWQC in no more than 20% of the cases (Suter and Tsao, 1996). Thus, this may result in an underestimation of risk.

A surface water benchmark was not available for 1,1,2-trichloro-1,2,2-

trifluoromethane. This VOC had the highest concentration of any of the VOCs

measured in groundwater, which may result in an underestimation of risk.

10.4.2 Risk Description

A coreening-level ecological risk assessment was conducted at this site to evaluate which

contaminants pose a potential adverse impact to ecological receptors inhabiting the Evergreen Manor

site and the adjacent Rock River and Dry Creek. Direct impacts on fish were evaluated for VOCs

discharging from groundwater to surface water. Actual measured concentrations in surface water

did not exceed detection limits, and the constituents detected in sediment were from upgradient

samples, or do not have available toxicological data. Thus, surface water and sediment pathways

were not evaluated.

To ensure that sites which may pose an ecological risk are properly identified, U.S. EPA (1996d)

suggests that values used in a screening level assessment should be consistently biased in the

direction of overestimating risk." Without this bias, a screening evaluation could not provide a

defensible conclusion for an absence of ecological risk."

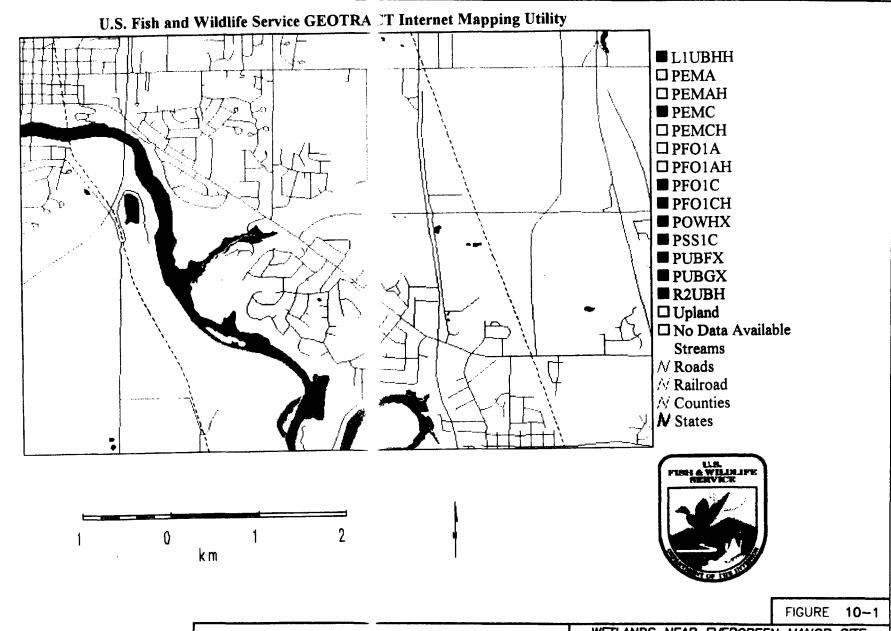
CHLAN01\WP\WO\RAC\036\29672-S10.WPD

RFW036-2A-AHVH

Section: 10 Revision: 1

Date: 28 March 2001 Page: 10 of 13

The screening-level ecological risk assessment found that there is a negligible potential for adverse effects on aquatic organisms in the Rock River and Dry Creek from site-related chemical contamination.



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EVERGREEN MANOR SITE Roscoe, Illinois

Section: 6 Revision: 1

Date: 28 March 2001 Page: 12 of 13

Table 10-1

# Preliminary E ological Conceptual Site Model Evergreen Manor Site Roscoe, Vinnebago County, Illinois

Exposure Medium	Exposure Route	Birds	Mammals	Fish	Macrobenthos
Groundwater	Ingestion	0	0	х	0
(discharge to surface water)	Dermal contact	0	0	х	0
	Inhalation	0	0	х	0
	Prey	0	O	х	О,

X = Potential exposure route determined to be significant for this receptor.

O = Potential exposure route determined to be insignificant or cannot be evalua d for this receptor.

Evergreen Manor Remedial Investigation Report Section: 10 Revision: 1 Date: 28 March 2001

Page: 13 of 13

Table 10-2

Comparison of Groundwate: Concentrations to Surface Water Screening Values

Evergreen Manor Site

Rosce, Winnebago County, Illinois

A I concentrations in ug/L

	Frequency of	Range of	Detec	วิทธ	Maximum	Tier I	Value	Lowest Chronic Value
Chemical	Detection	Minimum	Max	um	Location	Secondary acute	Secondary chronic	chronic
Residential wells	<del></del>							
cis-1,2-Dichloroethene	1/20		2.0	ō	RW-04	1100	590.00	9538*
Trichloroethene	3/20	0.7	6.	<u> </u>	RW-04	440	47.00	7257°
Tetrachloroethene	2/20	0.9	2.	$\overline{}$	RW-04	830	98.00	750
Acetone	2/20	0.6	0.	ĵ_	RW-03	28000	1500.00	507640°
1,1,1-Trichloroethane	6/20	0.6	5.	<u> </u>	RW-07	200	11.00	3493*
Chloroform	1/20	-	0.	5	RW-08	490	28.00	1240
Toluene	9/20	1	2.	2	RW-13;14;15;17;18;22	120	9.80	1269°
CPT wells								
Acetone	31/73	1	100	<u> </u>	CPT-05-06	28000	1500.00	507640*
1,1-Dichloroethane	1/73		2.	<u> </u>	CPT11-05		47.00	
cis-1,2-Dichloroethene	5/73	1	2.	<u> </u>	CPT-01-02	1100	590.00	9538*
Trichloroethene	8/73	2	4.	ெ	CPT-01-03;01-04;01-06	440	47.00	7257*
Toluene	60/73	0.5	3.	<u> </u>	CPT-11-08	120	9.80	1269*
Methylene chloride	1/73		0.	<u> </u>	CPT-03-05	26000	2200.00	42667
Benzene	3/73	0.5	0.	0	CPT-09-07	2300	130.00	525000
2-Butanone	11/73	-	16	ō	CPT-05-06	240000	14000.00	282170°
m-&/or p-Xylene	3/73	0.5	0.	5	CPT-06-01;09-07	32	1.80	62308° a
Xylenes (total)	2/73	0.6	0.	<u> 5 </u>	CPT11-01;20-01	32	1.80	62308° a
Ethylbenzene	1/73	-	0	5	CPT-09-07	130	7.30	>440
Tetrachloroethene	2/73	0.6	0.	5	CPT-10-02	830	98.00	750
1,1,1-Trichloroethane	17/73	0.8	3.	5	CPT-11-06	200	11.00	3493*
Monitoring wells								
Acetone	1/15	-	35	ō	MW103D	28000	1500.00	507640°
cis-1,2-Dichloroethene	2/15	1	2.	$\overline{\mathfrak{I}}$	MW105D	1100	590.00	9538*
Trichloroethene	3/15	2	3.	ĵ	MW105D	440	47.00	7257°
Tetrachloroethene	1/15	1	9.	ĵ.	MW103S	830	98.00	750
1,1,2-Trichloro-1,2,2-trifluoromethane	2/15	1	300	<b>90</b>	MW103D	NA NA	NA NA	NA
1,1,1,-Trichloroethene	6/15	1	3	Ū	MW103D:105D	200	11.00	3493°

#### Notes:

Only chemicals measured above detection limits are presented.

No Illinois general use water quality standards or national ambient water quality criteria available for the COPCs.

Tier II and lowest chronic values from Suter and Tsao, 1996.

NA = No available.

\* = Estimated value (Suter and Tsao, 1996).

a = Value is for total xylenes (Suter and Tsao, 1996).

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RFW036-2A-AHVH

Evergreen Manor

Remedial Investigation Report

Section: 11 Revision: 1

Date: 28 March 2001

Page: 1 of 7

**SECTION 11** 

CONCLUSIONS AND RECOMMENDATIONS

This section presents conclusions and recommendations for actions at the Evergreen Manor site,

based on the results of this remedial investigation. The conclusions subsection summarizes the major

findings of this report, and the recommendations subsection presents recommendations for future

action.

11.1 CONCLUSIONS

The following subsections summarize the major findings of this report with regard to site geology

and hydrogeology, nature and extent of contamination, contaminant fate and transport, and the

assessment of risks presented by the contaminated media.

11.1.1 Site Geology and Hydrogeology

The geology of the site is characterized by fill, topsoil, and/or organic deposits overlying thick

glaciofluvial outwash deposits of sand and gravel. These deposits in-filled the former Rock River

Valley during and following the most recent glacial events. Although none of the CPT

investigations reached bedrock, previous studies (Berg, et.al., 1981; Wehrmann, 1984) indicated that

overburden deposits in the vicinity of the Evergreen Manor site are approximately 250 feet thick and

overlie the bedrock surface. The hydraulic conductivity of this unconfined aquifer has been

estimated by others to be 3.8x10<sup>-2</sup> cm/sec (Wehrmann, 1984).

Based on measurements collected during the RI, the depth to groundwater was found to vary

between approximately 30 to 39 feet below grade. Table 6-1 indicates that depth to water in well

cluster MW101 is about 3 ft bgs, however, this well cluster is located in the Roscoe Sand and Rock

CHLAN01\WP\WO\RAC\036\29672S-11.WPD

RFW036-2A-AHVH

Section: 11 Revision: 1

Date: 28 March 2001

Page: 2 of 7

quarry, approximately 30 ft lower in elevation than McCurry Road. Similarly, well cluster 110 is

located in a depression along Illinois Route 251.

When measured in well clusters, groundwater elevations were found to be nearly identical. This

indicates the shallow and deeper aquifer zones are in direct hydraulic connection and that there are

no zones of fine material (i.e., silt or clay) that would create semi-confining zones. Based on

stratigraphy data from the CPT investigation, the aquifer is believed to be unconfined, and is most

likely continuous down to bedrock. The horizontal hydraulic gradient, based on groundwater

elevations shown on Figure 6-5, is approximately 0.0015 ft/ft. Using this gradient, the hydraulic

conductivity of 3.8x10<sup>-2</sup> cm/sec, and an estimated porosity of 30% (Fetter, 1994), the average linear

flow velocity was calculated to be 1.9x10<sup>-4</sup> cm/sec (0.54 ft/day).

11.1.2 Nature and Extent of Contamination

Screening levels were developed for illustrating the nature and extent of contamination at this site

and to identify COCs in antitronmental media at the site. The constituents detected in each medium

were compared to the screening levels. Concentrations of constituents exceeding the screening

levels were considered to be COCs. Screening levels are not directly implemented as RAOs; these

may be developed, as appropriate, in the future FS/RA process. The following paragraphs identify

the COCs found, and the nature and extent of contamination.

Groundwater

A total or 13 VOCs, as shown on Tables 7-2 through 7-4, were detected in samples collected from

CPT, monitoring wells, and residential wells. Only three of these constituents exceeded screening

levels. Trichloroethene was detected in residential well RW04 (11990 Blue Spruce Dr.), at a

concentration of 6 ppb, which exceeded the screening level of 5 ppb. Chloroform was detected in

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RFW036-2A-AHVH

Evergreen Manor

Remedial Investigation Report

Section: 11 Revisión: 1

Date: 28 March 2001

Page: 3 of 7

residential well RW08 (11943 Wagon Ln.), at a concentration of 0.9 ppb, which exceeded the

screening level of 0.02 ppb. Tetrachloroethene was detected in monitoring well MW103S at a

concentration of 9 ppb, which exceeded the screening level of 5 ppb. The levels of chlorinated

VOCs detected during this RI are significantly lower than those previously determined, indicating

continued diminishing contamination of the aquifer.

Based on these detections, chloroform, tetrachloroethene, and trichloroethene are considered COCs.

The extent of positive detections is shown on Figure 7-1, however, the extent of contamination

should only be considered as the area in the immediate vicinity of the three locations where COCs

were identified.

**Sediments** 

A total of four VOCs, as shown in Table 7-5, were detected in the collected sediment samples from

Dry Creek and the Rock River. The detected concentrations of the constituents were orders of

magnitude loss than their respective cornaning lawels. Therefore there are no COCs in the codiment

medium.

Surface Water

As shown on Table 5-8, the results of the surface water analyses were all below the method detection

limits. Therefore, there are no COCs in the surface water medium.

Sources

Previous investigations have not been able to pinpoint an exact source of contamination for the

Evergreen Manor site, but have identified the industrial area near the intersection of Rockton Road

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RFW036-2A-AHVH

Section: 11 Revision: 1

Date: 28 March 2001

Page: 4 of 7

and Illinois Route 251 as the likely source area of contamination. The same general conclusion is

drawn from the results of this RI. This conclusion is based on the results of the groundwater

analytical data from this RI and the fracture trace analysis results. Concentrations detected were not

able to lead to the identification of an exact source, but the fracture trace analysis indicated that the

industrial area near Rockton Road and Illinois Route 251 is a potential source area.

11.1.3 Contaminant Fate and Transport

It is possible for COCs identified at the site to migrate from one medium to another. COCs at the

site were only found in groundwater. Through groundwater movement these COCs migrate

downgradient by advection and discharge into the Rock River. The concentrations of COCs will

decrease during transport due mostly to dispersion, and possibly to biodegradation. Thus, when the

COCs enter the Rock River, and are diluted, concentrations of COCs are below analytical detection

limits. Although undetectable in surface water, the groundwater to surface water pathway is

considered a pathway of concern. Other pathways, such as the groundwater to air pathway, are not

considered to make a threat

The BIOSCREEN model was used to estimate when COC concentrations will decline below

screening levels. The model results predict that chloroform will be below its screening level about

three years after the RI, in about 2003. TCE is predicted to be below its screening level about six

years after the RI, in about 2006. PCE is predicted to be below its screening level about 15 years

after the RI, in about 2015.

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RFW036-2A-AHVH

Section: 11 Revision: 1

Date: 28 March 2001

Page: 5 of 7

11.1.4 Risk Assessment

Baseline human health and ecological risk assessments were performed to evaluate the potential

human health and environmental impacts associated with the site if no removal actions were

undertaken. The following subsections summarize the results of each risk assessment.

11.1.4.1 Baseline Human Health Risk Assessment

The baseline human health risk assessment evaluated residential users, and commercial/industrial

workers as the primary receptors at the site. The maximum detected concentration was used as the

exposure point concentration combining data from the residential wells, monitoring wells, and CPT

since no identifiable plume was found. Potential exposure to site groundwater was estimated

individually for adult and child residents and adult commercial/industrial workers. An exposure

scenario with a calculated cancer risk value exceeding 1E-06, and/or a non-cancer hazard index

exceeding 1.0, is considered to have risk.

The residential groundwater exposure scenario carcinogenic risk estimates ranged from 4.6E-06 to

1.9E-05. The constituents that primarily drove the carcinogenic risk numbers in these scenarios

include chloroform, tetrachloroethene, and trichloroethene. The residential groundwater exposure

scenario total hazard index ranged from 1.4 to 1.4 for an adult and from 3.7 to 3.8 for a child. The

risk was primarily driven by chloroform across the inhalation pathway for both children and adults.

To a lesser extent, trichloroethene contributed to the inhalation pathway risk value as well.

The commercial/industrial land use exposure scenario carcinogenic risk estimates ranged from

2.0E-06 to 6.9E-06. The risk was primarily driven by tetrachloroethene, trichloroethene, and

chloroform. The commercial/industrial land use exposure scenario total hazard index ranged from

0.97 to 0.99, which does not exceed the non-cancer threshold of unity.

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RFW036-2A-AHVH

Section: 11 Revision: 1

Date: 28 March 2001

Page: 6 of 7

11.1.4.2 Ecological Risk Assessment

The ecological risk assessment focused on the potential direct impact of VOCs to fish living in Dry

Creek and the Rock River. The screening-level ecological risk assessment found that there is

negligible potential for adverse effect on aquatic organisms in the Rock River and Dry Creek from

site-related chemical constituents.

11.1.5 Contaminant Concentrations

Based on a comparison of data presented in this RI and the data used to support the Hazard Ranking

Score, it appears that the concentrations of chlorinated VOCs are decreasing. There has been an

approximate order of magnitude decline in concentrations.

11.1.6 Non-Time Critical Removal Action

Recause concentrations of VOCs were still detected above screening levels, an appropriate remedy

is warranted under current guidance and protocol. However, the U.S. EPA completed a non-time

critical removal action at the Evergreen Manor site in September 2000. This included the installation

of an extension to the existing public water system to include residences located in the subdivisions

impacted by VOCs, and the abandonment of private residential water wells. This removal action

is discussed in greater detail in the EE/CA, and the Action Memorandum requesting its

implementation.

The result of this removal action is that it has effectively deleted the residential well exposure

pathway that was discussed in the human health risk assessment. Thus, since the exposure pathway

has been eliminated, the associated human health risk has also been eliminated.

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RFW036-2A-AHVH

Section: 11 Revisión: 1

Date: 28 March 2001

Page: 7 of 7

# 11.2 RECOMMENDATIONS

Based on the conclusions put forth above, there does not appear to be significant risk associated with the Evergreen Manor site at this time. The following recommendations are intended to make sure that no new risks could develop as a result of changing site conditions or new developments.

- Although risk to residents in the Evergreen Manor site subdivisions has been addressed, not all of the residences were included in the public water system expansion. Residential wells still used to provide water from the shallow aquifer to residences on either side of the expanded water system should be sampled for VOCs at least every five years. This will ensure that the removal action has met its goals and will protect residents close to the contamination.
- Based on a historical decline of concentrations across the site, as evidenced by laboratory data, and as simulated through modeling, it is expected that constituent concentrations will continue to decline, and ultimately decline below screening levels. The IEPA-installed monitoring wells should be sampled periodically in order to ensure that constituent concentrations are indeed declining, and that ultimately no more contamination travels to the remaining residential wells.
  - or near the Evergreen Manor site contamination. This could be accomplished through local ordinance or deed restrictions and should cover those areas not currently used for residential purposes.
- As a result of the RI work performed, no further attempts at source identification are recommended. Thus, soil and sediment sampling is not warranted, and no new monitoring wells are recommended at this time.
- As required by the U.S. EPA, a Five Year Review of the Evergreen Manor site should be completed. This review should evaluate any groundwater data collected since this RI, and evaluate the effectiveness, with respect to areal coverage, of the public water supply in the Evergreen Manor site subdivisions and other areas coincident with the plume.

Evergreen Manor Remedial Investigation Report Section: 12 Revision: 1

Date: 28 March 2001 Page: 1 of 4

### **SECTION 12**

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Section: 12 Revision: 1 Date: 28 March 2001

Page: 2 of 4

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Evergreen Manor Remedial Investigation Report Section: 12 Revision: 1

Date: 28 March 2001 Page: 4 of 4

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# APPENDIX A

# SUPPORTING MATERIAL FOR HUMAN HEALTH AND ECOLOGICAL RISK ASSESSMENTS





# United States Department of the Interior

### FISH AND WILDLIFE SERVICE

Ecological Services
Rock Island Field Office
4469 48th Avenue Court
Rock Island, Illinois 61201
Tel: 309/793-5800 Fax: 309/793-5804

July 10, 2000

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Ce K Fischer T. Bosko

Mr. Terry Bosko Technical Manager Roy F. Weston, Inc. 750 E. Bunker Ct. Ste 500 Vernon Hills, IL 60061-1450

Dear Mr. Bosko:

This is in response to your letter to our Rock Island Field Office dated June 27, 2000, requesting threatened and endangered species information relative to your Evergreen Manor project area. Our comments are provided below.

We have reviewed the information provided for the location of this site. To facilitate compliance with Section 7(c) of the Endangered Species Act of 1973, as amended, Federal agencies are required to stain from the Fish and William State information personnel and species, listed or proposed to be listed, which may be present in the area of a proposed action. Therefore, we are furnishing you the following list of species which may be present in the

concerned area:

Classification	Common Name (Scientific Name)	<u>Habitat</u>
Endangered	Indiana bat (Myotis sodalis)	Caves, mines; small stream corridors with well developed riparian woods; upland forests
Threatened	Prairie bush clover (Lespedeza leptostachya)	Dry to mesic prairies with gravelly soil
Threatened	Bald Eagle	Wintering, breeding

This is your future. Don't leave it blank. - Support the 2000 Census.

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The endangered Indiana bat (Myotis sodalis) could potentially occur throughout the state of Illinois. During the summer, the Indiana bat frequents the corridors of small streams with well developed riparian woods as well as mature upland forests. It forages for insects by flying beneath the tree canopy, and roosts and rears its young beneath the loose bark of large dead or dying trees. It winters in caves and abandoned mines.

Suitable summer habitat in Iowa is considered to have the following characteristics within a ½ mile radius of the project site:

- 1) forest cover of 15% or greater;
- 2) permanent water;
- 3) one or more of the following tree species 9 inches diameter at breast height (dbh) or greater: shagbark hickory, bitternut hickory, American elm, slippery elm, eastern cottonwood, silver maple, white oak, red oak, post oak, and shingle oak;
- 4) at least 1 potential roost tree per 2.5 acres;
- 5) potential roost trees must have greater than 10% coverage of loose bark (by visual estimation of peeling bark on trunks and main limbs).

If the project site contains any habitat that fits the above description, it may be necessary to conduct a survey to determine whether the bat is present. If Indiana bats are known to be present, they must not be harmed, harassed, or disturbed and their habitat must not be destroyed. Indiana bat habitat may be altered only between the dates of October 1 and March 31.

The prairie bush alover (Leanedeze Intestacliva) is listed as threatened in Winnehoge County in Illinois. It occupies dry to mesic prairies with gravelly soil. There is no critical habitat designated for this species. Federal regulations prohibit any commercial activity involving this species or the destruction, malicious damage or removal of this species from Federal land or any other lands in knowing violation of State law or regulation, including State criminal trespass law. This species should be searched for whenever prairie remnants are encountered.

The threatened bald eagle (Haliaeetus leucocephalus) is listed as breeding and wintering in Winnebago County. During the winter, this species feeds on fish in the open water areas created by dam tailwaters, the warm water effluents of power plants and municipal and industrial discharges, or in power plant cooling ponds. The more severe the winter, the greater the ice coverage and the more concentrated the eagles become. They roost at night in groups in large trees adjacent to the river in areas that are protected from the harsh winter elements. They perch in large shoreline trees to rest or feed on fish. There is no critical habitat designated for this species. The eagle may not be harassed, harmed, or disturbed when present nor may nest trees be cleared.

These comments are provided under the authority of and in accordance with the provisions of the Fish and Wildlife Coordination Act (16 U.S.C. 661 et seq.; 48 Stat. 401), as amended; and the Endangered Species Act of 1973, as amended.

Mr. Terry Bosko 3.

Thank you for the opportunity to provide comments early in the planning process. If you have any additional questions or concerns, please contact Kevin de la Bruere of my staff at extension 530.

Sincerely

Supervisor

# TABLE I SEL CTION OF EXPOSURE PATHWAYS EVERGREEN MANC & SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

	<del></del>	w=	T	T			T	<del></del>	
	Medium	Evangues		Personal Control	Panents-	E	On Sind	Turnes	Batianala for Calentina or Evaluation
Scenario	Medium	Exposure	Exposure	Receptor	Receptor	Exposure	On-Site/	Type of	Rationale for Selection or Exclusion
Timeframe		Medium	Point	Population	Age	Route	Off-Site	Analysis	of Exposure Pathway
A	Residential well	Groundwater	Tap water	Resident					
Current	Residential Well	Oroundwater	Tap water	Kesiden	Adult	Ingestion	On-site	Quant	Private wells currently in use.
						Dermal	On-site	Quant	Private wells currently in use
	[ ]				Child	Ingestion	On-site	Quant	Private wells currently in use
	ļ į					Dermai	On-site	Quant	Private wells currently in use
		Air	Water vapors at showerhead	Resident	Adult	Inhalation	On-site	Quant	Private wells currently in use.
					Child	Inhalation	On-site	Quant	Private wells currently in use.
Future	Monitoring well	Groundwater	Tap water	Resident	Adult	Ingestion	On-site	Quant	Private well may potentially be installed in the future
						Dermal	On-site	Quant	Private well may potentially be installed in the future.
				·	Child	Ingestion	On-site	Quant	Private well may potentially be installed in the future.
						Dermal	On-site	Quant	Private well may potentially be installed in the future.
		Air	Water vapors at showerhead	Resident	Adult	Inhalation	On-site	Quant	Private well may potentially be installed in the future.
	İ			Í	Child	Inhalation	On-site	Quant	Private well may potentially be installed in the future.
Future	Monitoring well	Groundwater	Tap water	Commercial/Industria	Adult	Ingestion	On-site	Quant	Private well may potentially be installed in the future.
						Dermal	On-site	Quant	Private well may potentially be installed in the future.
		Air	Water vapors at showerhead	Commercial/Industria	Adult	Inhalation	On-site	Quant	Private well may potentially be installed in the future.
Future	CPT well	Groundwater	Tap water	Resident	Adult	ingestion	On-site	Quant	Private well may potentially be installed in the future
				•		Dermal	On-site	Quant	Private well may potentially be installed in the future
					Child	Ingestion	On-site	Quant	Private well may potentially be installed in the future
						Dermal	On-site	Quant	Private well may potentially be installed in the future
		Air	Water vapors at showerhead	Resident	Adult	Inhelation	On-site	Quant	Private well may potentially be installed in the future
					Child	Inhelation	On-site	Quant	Private well may potentially be installed in the future
uture	CPT well	Groundwater	Tap water	Commercial/Industria	Adult	Ingestion	On-site	Quant	Private well may potentially be installed in the future
			1			Dermal	On-site	Quant	Private well may potentially be installed in the future
		Air	Water vapors at showerhead	Commercial/Industria:	Adult	Inhalation	On-site	Quant	Private well may potentially be installed in the future

#### TABLE 2.1

## OCCURRENCE, DISTRIBUTIO: AND SELECTION OF CHEMICALS OF POTENTIAL CONCERN EVERGREEN MANOR SITE

ROSC DE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Current/Future

Medium: Groundwater

Exposure Medium: Groundwater

Exposure Point: Tap water

CAS	Chemical	Minimum	Minimum	Maximum	Maximum	Units	Location	Detection	Range of	Concentration	Background	Screening	Potential	Potential	COPC	Rationale for (4)
Number		Concentration	Qualifier	Concentration	Qualifier		of Maximum	Frequency	Detection	Used for	Value	Toxicity Valu	ARAR/TBC	ARAR/TBC	Flag	Contaminant
							Concentration		Limits	Screening (2)			Value	Source		Deletion or Selection
156-59-2	cis-1,2-Dichloroethene	1	J	2	J	ug/L	WW106S;106D;CPT-01-0	8/108	4	2	N/A	6.1	70	IEPA-TACO	N	BSL
75-34-3	1,1-Dichloroethane		_	2.0	J	ug/L	CPT11-05	1/108	1	2.0	N/A	70	700	IEPA-TACO	N	BSL
79-01-6	Trichloroethene	0.7	J	6	-	ug/L	RW-04	14/108	1	6	N/A	0.16	5	IEPA-TACO	Y	ASL
127-18-4	Tetrachloroethene	0.6	j	9	J	ug/L	MW103S	5/108	1	9	N/A	0.11	5	IEPA-TACO	Y	ASL
67-64-1	Acetone	0.6	_	100.0	J	ug/L	CPT-05-08	34/108	2	100.0	N/A	61	700	IEPA-TACO	N	ASL
71-55-6	1,1,1-Trichloroethane	0.6	j	5	J	ug/L	RW-07	29/108	2	5	N/A	79	200	IEPA-TACO	N	BSL
	1,1,2-Trichloro-1,2,2-trifluoromethane	2	J	300	J	ug/L	MW103D	2/108	2	300	N/A	N/A	N/A	IEPA-TACO	Z	NTX
67-66-3	Chloroform	-	ļ	0.9	J	ug/L	RW-08	1/108	3	0.9	N/A	0.02	0.02	IEPA-TACO	Υ	ASL
108-88-3	Toluene	1	j	3	-	ug/L	CPT-11-08	69/108	1	3	N/A	72	1000	IEPA-TACO	N	BSL
75-09-2	Methylene chloride	-	-	0.5	J	ug/L	CPT-03-05	1/108	1	0.5	N/A	0.43	5	IEPA-TACO	Υ	ASL
71-43-2	Benzene	0.5	J	0.6	J	ug/L	CPT-09-07	3/108	1	0.6	N/A	0.04	5	IEPA-TACO	Υ	ASL
78-93-3	2-Butanone	-	-	16.0	J	υg/L	C PT-05-06	1/108	11	16.0	N/A	190.00	N/A	N/A	N	BSL
1330-20-7	Xylenes (total)	0.6	J	0.6	J	ug/L	CPT11-01;02-01	2/108	1	0.6	N/A	140.00	10000	IEPA-TACO	N	BSL
1330-20-7	m-&/or p-Xylene	0.5	J	0.7	J	ug/L	CPT-06-01;02-07	3/108	1	0.7	N/A	140.00	10000	IEPA-TACO	Ν	BSL
100-41-4	Ethylbenzene	-		0.6	J	ug/L	CPT-09-07	1/108	1	0.6	NA	130.00	700	IEPA-TACO	2	BSL

- (1) Minimum/maximum detected concentration.
- (2) Maximum detected concentration used as screening value
- (3) Background value not available for groundwater.
- (4) IEPA-TACO (35 IAC 740). See Section 2 for supporting information. For screening toxicity values: Cancer benchmark value = 1-07; HQ = 0.1.
- (5) Rationale Codes Selection Reason:

Infrequent Detection but Associated Historically (HIST)

Frequent Detection (FD)

Toxicity Information Available (TX)

Above Screening Levels (ASL)

Deletion Reason:

Infrequent Detection (IFD)

Background Levels (BKG)

No Toxicity Information (NTX)

Essential Nutrient (NUT)

Below Screening Level (BSL)

Definitions: N/A = Not Applicable

SQL = Sample Quantitation Limit

COPC = Chemical of Potential Concern

ARAR/TBC = Applicable or Relevant and Appropriate Requirement/To Be Considered

UMTRCA = Uramium Mill Tailings Radiation Control Act soil protection standard (40 CFR 192).

MCL = Maximum contaminant level.

J = Estimated Value

C = Value has been changed from original data sheet

X = Gross value (no instrument background subtracted); actual net value will be approximately 1.2 pC

TABLE 3.1

MEDIUM-SPECIFIC EX 'OSURE POINT CONCENTRATION SUMMARY EVERGREEN MANOR S (E, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Cufrent/Future

Medium: Groundwater

Exposure Medium: Groundwater

Exposure Point: Tap water

Chemical of	Units	Arithmetic Mean	95% UCL of Normal	Maximum Detected	i	imum alifler	EPC Units	1	easonable Ma	ximum Exposure		Centr	al Tendency
Potential			Data	Concentration				Medium	Medium	Medium	Medium	Medium	Medium
Concern								EPC	EPC	EPC	EPC	EPC	EPC
					L			Value	Statistic	Rationale	Value	Statistic	Rationale
Trichloroethene	ug/L			6		J	ug/L	6	MAX	Undefined center of plume	6	MAX	Undefined center of plume
Tetrachloroethene	ug/L			9		j	ug/L	9	MAX	Undefined center of plume	9	MAX	Undefined center of plume
Chloroform	ug/L	-		0.9		J	ug/L	0.9	MAX	Undefined center of plume	0.9	MAX	Undefined center of plume
Acetone	ug/L	-		100		J	ug/L	100	MAX	Undefined center of plume	100	MAX	Undefined center of plume
Benzene				0.6		J	ug/L	0.6	MAX	Undefined center of plume	0.6	MAX	Undefined center of plume
Methylene chloride	ug/L	<b></b>		0.5		J	ug/L	0.5	MAX	Undefined center of plume	0.5	MAX	Undefined center of plume

Statistics: Maximum Detected Value (Max); 95% UCL of Normal Data (95% UCL-N); 95% UCL of Log-transformed Data (95% UCL-T); Mean of Log-transformed Data (Mean-Mean of Normal Data (Mean-N).

#### TABLE 4.1a

#### VALUES I SED FOR DAILY INTAKE CALCULATIONS

EVERGREEN MAN-IR SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Current/Future
Medium: Groundwater
Exposure Medium: Groundwater
Exposure Point: Tap water
Receptor Population: Resident
Receptor Age: Adult

Exposure Route	Parameter Code	Parameter Definition	{	its	RME Value	RME Rationale/ Reference	CT Value	CT Rationale/ Reference	Intake Equation/ Model Name
ngestion	ĊW	Chemical concentration in water		χĹ	See Table 3	See Table 3	See Table 3	See Table 3	
	IR-W	Ingestion rate of water	) i	iay	2	EPA, 1999	1.4	EPA, 1997	CDI (mg/kg/d) = CW x IR x 1E-03
	EF	Exposure frequency	d.	s/yr	350	EPA, 1999	350	EPA, 1991c	mg/ug x EF x ED / (BW x AT)
	EDc	Exposure duration -carcinogens		r <b>s</b>	24	EPA, 1999	7	EPA, 1997	
	EDnc	Exposure duration - noncarcinogens		<b>(\$</b>	30	EPA, 1999	9	EPA, 1997	
	8W	Body weight		9	70	EPA, 1999	70	EPA, 1999	
	AT-C	Averaging time (cancer)		ув	25550	EPA, 1999	25550	EPA, 1999	
	AT-NC	Averaging time (non-cencer)		ys.	8760	EPA, 1999	2555	EPA, 1999	
Dermal absorption	cw	Chemical concentration in water	<del>                                     </del>	<u>,,</u>	See Table 3	See Table 3	See Table 3	See Table 3	<b></b>
	SA	Surface area available for contact		n²	23000	EPA, 1992a	20000	EPA, 1992a	
	PC	Permeability constant		√hr	chemical specific	EPA, 1992a	chemical specific	EPA, 1992a	CDI (mg/tg/d) = CW x SA x PC 1L/1000 cm3 x ET x EF x ED x 1/BW x 1/AT
	ET	Exposure time	}	day	0.75	EPA, 1997	0 33	EPA, 1997	
	EF	Exposure frequency	1	s/vr	350	EPA, 1991c	350	EPA, 1991c	
	EDc	Exposure duration -carcinogens		3	24	EPA. 1999	7	EPA, 1997	
	EOnc	Exposure duration - noncarcinogens	1		30	EPA, 1999		EPA, 1997	
	BW	Body weight		g	70	EPA, 1999	70	EPA, 1999	
	AT-C	Averaging time (cancer)		.ys	25550	EPA, 1999	25550	EPA, 1999	
	AT-NC	Averaging time (non-cancer)		:ys	8760	EPA, 1999	2555	EPA, 1999	
nhalation of vapors	cw	Chemical concentration in water		уL	See Table 3	See Table 3	See Table 3	See Table 3	
	IR-A	Inhalation rate of indoor air	-	day	15	EPA, 1999	15	EPA, 1999	
	EF	Exposure frequency	e	/NT	350	EPA, 1999	350	EPA, 1999	CDI (mg/kg-d) = CW x IR-A x E ED x K x 1/BW x 1/AT
	EDc	Exposure duration -carcinogens		r <b>s</b>	24	EPA, 1999	7	EPA, 1997	
	EDnc	Exposure duration - noncarcinogens		rs.	30	EPA, 1999	9	EPA, 1997	
ļ	BW	Body weight	1	9	70	EPA, 1999	70	EPA, 1999	
	AT-C	Averaging time (cancer)	.	ıys	25550	EPA, 1999	25550	EPA, 1999	
	AT-NC	Averaging time (non-cancer)		.ys	8760	EPA, 1999	2555	EPA, 1999	
									1
ļ	VF	Volaitization factor	. u.	1005	5,00E-01	EPA 1999	5.00E-01	EPA 1999	ļ

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#### TABLE 4.1b

# VALUES SED FOR DAILY INTAKE CALCULATIONS EVERGREEN MAN. R. SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Curren/Future
Medium: Groundwater
Exposure Medium: Groundwater
Exposure Point: Tap water
Receptor Population: Resident
Receptor Age: Child

Exposure Route	Parameter Code	Parameter Definition	Uts	RME Value	RME Rationale/ Reference	CT Value	CT Rationale/ Reference	intake Equation/ Model Name
Ingestion	cw	Chemical concentration in water	U L	See Table 3	See Table 3	See Yable 3	See Table 3	
l j	IR-W	Ingestion rate of water	Lay	1	EPA, 1999	0.7	EPA, 1997	CDI (mg/kg/d) = CW x IR x 1E-03
	EF	Exposure frequency	da <sub>3</sub> ./yr	350	EPA, 1999	350	EPA, 1999 Professional judgement;	mg/ug x EF x ED / (BW x AT)
	ED	Exposure duration	٠,٠	6	EPA, 1999	2	see Sec. 3.3	
	8W	Body weight	٠,	15	EPA, 1999	15	EPA, 1999	
	AT-C	Averaging time (cancer)	d. s	25550	EPA, 1999	25550	EPA, 1999	
	AT-NC	Averaging time (non-cancer)	d. s	2190	EPA, 1999	730	EPA, 1999	
Dermal absorption	cw	Chemical concentration in water	m i	See Table 3	See Table 3	See Table 3	See Table 3	
:	SA	Surface area available for contact	c '	8538	EPA., 1997	7314	EPA, 1997	
	PC	Permeability constant	cr: hr	chemical specific	EPA, 1992a	chemical specific	EPA, 1992a	CDI (mg/kg/d) = CW x SA x PC x 1L/1000 cm3 x ET x EF x ED x 1/BW x 1/AT
	ET	Exposure time	hrs any	0.75	EPA, 1997	0.33	EPA, 1997	
	EF	Exposure frequency	da <sub>)</sub> /yr	350	EPA, 1999	350	EPA, 1999 Professional judgement;	
	ED	Exposure duration	٠,٠	6	EPA, 1999	2	see Sec. 3.3	
	BW	Body weight	;	15	EPA, 1999	15	EPA, 1999	
	AT-C	Averaging time (cancer)	d s	25560	EPA, 1999	25550	EPA, 1999	
	AT-NC	Averaging time (non-cancer)	di. 5	2190	EPA, 1999	730	EPA, 1999	
nhalation of vapors	cw	Chemical concentration in water	m L	See Table 3	See Table 3	See Table 3	See Table 3	
	IR-A	Inhalation rate of indoor sir	m₃ ayr	8.7	EPA, 1999	8.7	EPA, 1999	
	EF	Exposure frequency	da yr	350	EPA, 1999	350		CDI(mg/kg-d) = CW x IR-A x EF x ED x K x 1/8W x 1/AT
	ED	Exposure duration	у,	6	EPA, 1999	2	see Sec. 3.3	
	BW	Body weight		15	EPA, 1999	15	EPA, 1999	
	AT-C	Averaging time (cancer)	Clas S	25550	EPA, 1999	25550	EPA, 1999	
	AT-NC	Averaging time (non-cancer)	di s	2190	EPA, 1999	730	EPA, 1999	
	VF	Volailization factor	uni: :se	5.00E-01	EPA 1999	5.00E-01	EPA, 1999	•

# TABLE 4.2 VALUES SED FOR DAILY INTAKE CALCULATIONS EVERGREEN MAN . R SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Current/Future

Medium: Groundwater

Exposure Medium: Groundwater

Exposure Point: Tap water

Receptor Population: Commercial/Industrial

Receptor Age: Worker

Exposure Route	Parameter Code	Parameter Definition		nits	RME Value	RME Rationals/ Raterence	CT Value	CT Rationale/ Reference	Intake Equation/ Model Name
Ingestion	cw	Chemical concentration in water		o/L	See Table 3	See Table 3	See Yable 3	See Table 3	
	IR-W	Ingestion rate of water		day	1	EPA, 1999	0.7	EPA, 1997	CDI (mg/kg/d) = CW x IR x 1E-03
	EF	Exposure frequency	c	/s/yr	250	EPA, 1999	250	EPA, 1991c	mg/ug x EF x ED / (BW x AT)
	EOc	Exposure duration -carcinogens		rs.	25	EPA, 1999	9	EPA, 1997	
	EDnc	Exposure duration - noncercinogens		.r <b>s</b>	25	EPA, 1999	9	EPA, 1997	
	BW	Body weight	ĺ	<b>₹</b> Ø	70	EPA, 1999	70	EPA, 1999	
	AT-C	Averaging time (cancer)		198	25550	EPA, 1999	25550	EPA, 1999	
	AT-NC	Averaging time (non-cancer)		1 <b>78</b>	9125	EPA, 1999	2555	EPA, 1999	
Dermal absorption	cw	Chemical concentration in water		₹.	See Table 3	See Table 3	See Table 3	See Table 3	
	SA	Surface area available for contact		m²	23000	EPA, 1992a	20000	EPA, 1992a	
					chemical specific	EPA, 1992a	chemical specific	EPA, 1992a	CDI (mg/kg/d) = CW x SA x PC x 1L/1000 cm3 x ET x EF x ED x 1/8W x 1/AT
	PC	Permeability constant		vhr	l '	EPA, 1997	•	EPA. 1997	1/201 A 1/A1
	ET	Exposure time	h	•	0.75		0.33	1	
	EF	Exposure frequency	С	/s/yr	250	EPA, 1991c	250	EPA, 1991c	
	EDc	Exposure duration -carcinogens		rs		EPA, 1999	9	EPA, 1997	
		Exposure duration - noncarcinogens		rs	25	EPA, 1999	9	EPA, 1997	
}		Body weight	l	.9	70	EPA, 1999	70	EPA, 1999	
	AT-C	Averaging time (cancer)		198		EPA, 1999	25550	EPA, 1999	
	AT-NC	Averaging time (non-cancer)		198	9125	EPA, 1999	2555	EPA, 1999	
nhelation of vapors	cw	Chemical concentration in water		g/L	See Table 3	See Table 3	See Table 3	See Table 3	
	IR-A	Inhalation rate of indoor air	r	day	15	EPA, 1999	15	EPA, 1999	CDI (mg/kg-d) = CW x IR-A x EF x
ŀ	EF	Exposure frequency	,	// <b>yr</b>	250	EPA, 1999	250	EPA, 1999	ED x K x 1/BW x 1/AT
	EDc	Exposure duration -carcinogens		r <b>s</b>	25	EPA, 1999	9	EPA, 1997	
	EDnc	Exposure duration - noncarcinogens		rs.	25	EPA, 1999	9	EPA, 1997	
	BW	Body weight		·g	70	EPA, 1999	70	EPA, 1999	
	AT-C	Averaging time (cancer)		ıys	25550	EPA 1999	25550	EPA, 1999	
	AT-NC	Averaging time (non-cancer)		1 <b>98</b>	9125	EPA, 1999	2555	EPA, 1999	
Ī	VF	Volaitization factor	U	Jess	5.00E-01	EPA, 1999	5.00E-01	EPA, 1999	l

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#### TABLE 5.1

#### NON-CA CER TOXICITY DATA - ORAL/DERMAL

#### EVERGREEN MANOR SITE

ROSC JE, WINNEBAGO COUNTY, ILLINOIS

								<u> </u>			
Chemical of Potential Concern	Chronic/ Subchronic	Oral RfD Value	Oral RfD Units	Oral to Adjustmen	ermal Factor (1)	Adjusted Dermal RfD (2)	Units	Primary Target Organ	Combined Uncertainty/Modifying Factors	Sources of RfD: Target Organ	Dates of RfD: Target Organ (3) (MM/DD/YY)
Acetone	Subchronic	1.00E-01	mg/kg-day			1.00E-01	mg/kg-day	liver/kidney	1000	IRIS	7/7/00
cis-1,2-Dichloroethene	Subchronic	1.00E-02	mg/kg-day			1.00E-02	mg/kg-day	blood	3000	HEAST/Region IX	11/29/99
Chloroform	Chronic	1.00E-02	mg/kg-day			1.00E-02	mg/kg-day	liver	1000	IRIS	7/7/00
Benzene	N/A	3.00E-03	mg/kg-day			3.00E-03	mg/kg-day	N/A	N/A	NCEA/Region IX	11/29/99
Trichloroethene	N/A	6.00E-03	mg/kg-day			6.00E-03	mg/kg-day	N/A	N/A	Withdrawn/Region IX	11/29/99
Toluene	Chronic	2.00E-01	mg/kg-day			2.00€-01	mg/kg-day	liver/kidney	1000	IRIS	7/7/00
Tetrachloroethene	Subchronic	1.00E-02	mg/kg-day			1.00E-02	mg/kg-day	liver	1000	IRIS	7/7/00
1,1,1-Trichtoroethane	N/A	3.50E-02	mg/kg-day	,		3.50E-02	mg/kg-day	N/A	N/A	NCEA/Region IX	11/29/99
1,1,2-Trichloro-1,2,2-trifluoromethane	N/A	N/A	N/A	N		N/A	N/A	N/A	N/A	N/A	N/A
Methylene chloride	Chronic	6.00E-02	mg/kg-day	١	:	6.00E-02	mg/kg-day	liver	100	IRIS	7/7 <b>/0</b> 0
2-Butanone	Chronic	6 00E-01	mg/kg-day			6.00E-01	mg/kg-day	decrease fetal body weight	3000	IRIS	7 <i>/71</i> 00
m-8/or p-Xylene	Chronic	2.00E+00	mg/kg-day			2.00E+00	mg/kg-day	hyperactivity, decreased body weight, increased mortality	100	IRIS	7 <i>1711</i> 00
Ethylbenzene	Subchronic	1.00E-01	mg/kg-day	,		1. <b>00</b> E-01	mg/kg-day	liver/kidney	1000	IRIS	7/7/00

#### N/A = Not Applicable

- (1) Refer to RAGS, Part A; Complete (100%) oral absorption is assumed due to a lack of scientifically defensible da . base on Gl absorption factors (EPA, 1989) See Section 4.
- (2) Dermai RfD assumed to equal oral RfD. See Section 4.
- (3) For IRIS values, provide the date IRIS was searched.

For HEAST values, provide the date of HEAST.

For NCEA values, provide the date of the article provided by NCEA.

Region IX = Region IX Risk-Based Concentration Table.

#### TABLE 5.2

# NON- ANCER TOXICITY DATA - INHALATION EVERGREEN MANOR SITE

RO : OE, WINNEBAGO COUNTY, ILLINOIS

Chemical	Chronic/	Value	Units	, ijusted	Units	Primary	Combined	Sources of	Dates (2)
of Potential	Subchronic	Inhalation	1	alation		Target	Uncertainty/Modifying	RIC:RID:	(MM/DD/YY)
Concern		RfC	ļ	rD (1)	•	Organ	Factors	Target Organ	(
2.1.2.11				]					
Acetone	Subchronic	N/A	N/A	0E-01	mg/kg-day	liver/kidney	1000	Route extrapolation Route extrapolation/Region	1/7/00
cis-1,2-Dichloroethene	Subchronic	N/A	N/A	.0E-02	mg/kg-day	blood	3000	ΙX	11/29/99
Chioroform	Chronic	3.00E-04	mg/m3	i 30 <b>E-05</b>	mg/kg-day	liver	1000	HEAST/Region IX	11/29/99
Benzene	N/A	6.00E-03	mg/m3	0E-03	mg/kg-day	N/A	N/A	NCEA/Region IX	11/29/99
Trichioroethene	N/A	2.10E-02	mg/m3	: >0E-03	mg/kg-day	N/A	N/A	Route extrapolation/Region	11/29/99
Toluene	Chronic	4.00E-01	mg/m3	0E-01	mg/kg-day	neurological	300	Route extrapolation	7/7/00
Tetrachiomethene	Subchronic	3.90E-01	mg/m3	10E-01	mg/kg-day	kver	1000	NCEA/Region IX	11/29/99
1,1,1-Trichloroethane	N/A	N/A	N/A	: +0E-01	mg/kg-day	N/A	N/A	NCEA/Region IX	11/29/99
1,1,2-Trichloro-1,2,2-trifluoromethane	N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A	N/A
Methylene chloride	N/A	N/A	N/A	->0E-01	mg/kg-day	N/A	N/A	HEAST/Region IX	11/29/99
2-Butanone	Chronic	1.00€+00	mg/m3	9E-1	mg/kg-day	decreased fetal body weight	3000	IRIS	7/700
m-&/or p-Xylene	N/A	N/A	N/A	. 0E-01	mg/kg-day	N/A	N/A	withdrawn/Region IX	11/29/99
Ethylbenzene	Chronic	1.00E+00	mg/m3	9E-1	mg/kg-day	developmental	300	IRIS	7/7/00
•			1			,			

#### N/A = Not Applicable

- (1) See Section 4 in text.
- (2) For IRIS values, provide the date IRIS was searched.

For HEAST values, provide the date of HEAST.

For NCEA values, provide the date of the article provided by NCEA.

Region IX = Region IX Risk-Based Concentration Table.

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TABLE 6.1

CANCER TO CICITY DATA - ORAL/DERMAL

EVI RGREEN MANOR SITE

ROSCOE, V. NNEBAGO COUNTY, ILLINOIS

Chemical of Potential Concern	Oral Cancer Slope Factor	Oral to Dermal Adjustment Factor	Adjusted Dermal C scer Slope Factor (1)	Units	Weight of Evidence/ Cancer Guideline Description	Source Target Organ	Date (2) (MM/DD/YY)
Acetone	N/A	N/A	NKA	N/A	D	IRIS	7/7/00
cis-1,2-Dichloroethene	N/A	N/A	NKA	N/A	D	IRIS	7/7/00
Chloroform	6.10E-03	1	6.10E-03	(mg/kg-day) <sup>-1</sup>	B2/all kidney tumors	IRIS	7/7/00
Benzene	1.5E-02 to 5.5E-02	1	1.5E-02 to 5.5E-02	(mg/kg-day) <sup>-1</sup>	A/leukemia	IRIS	7/7/00
Trichloroethene	1.10E-02	1	1.1 <b>0E-0</b> 2	(mg/kg-day) <sup>-1</sup>	N/A	NCEA/Region IX	11/29/99
Toluene	N/A	N/A	N/A	N/A	D	IRIS	7/7/00
Tetrachioroethene	5.20E-02	1	6.20E-02	(mg/kg-day) <sup>-1</sup>	liver	NCEA/Region IX	11/29/99
1,1,1-Trichloroethene	N/A	N/A	N/A	N/A	D	IRIS	7/7/00
1,1,2-Trichloro-1,2,2-trifluoromethane	NA	N/A	NA	N/A	N/A	N/A	N/A
Methylene chloride	7.50E-03	1	7.60E-03	(mg/kg-day) <sup>-1</sup>	B2/hepatocellular	IRIS	7/7 <b>/0</b> 0
2-Butanone	NA	N/A	N/A	N/A	D	IRIS	7/7/00
m-&/or p-Xylene	N/A	N/A	NA	N/A	D	IRIS	7/7/00
Ethylbenzene	N/A	N/A	NA	N/A	D	IRIS	7 <i>/71</i> 00

 Refer to RAGS Part A and Section 4 of text.
 Complete (1(100%) oral absorption is assumed due to a tack of scientifically defensible database on GI absorption factors (EPA, 1989).

(2) For IRIS values, provide the date IRIS was searched.

For HEAST values, provide the date of HEAST.

For NCEA values, provide the date of article provided by NCEA.

REGION IX = Region IX Preliminary remediation goals.

RPA = Relative Potency Approach (EPA, 1993a)

EPA ( rouge

- A -- Ruman carcinogen
- B1 Probable human carcinogen indicates that limited human data are available

- B2 Probable human carcinogen indicates sufficient evidence in animals and inadequate or no evidence in humans
- C Possible human carcinogen
- D Not dissillable as a human carcinogen
- E Evidence of noncercinogenicity

Weigt of Evidence:

Kn wryLibrary

Ca not be Determined

No Likely

#### TABLE 6.2

#### CAN :ER TOXICITY DATA -- INHALATION

#### **EVERGREEN MANOR SITE**

ROSC DE, WINNEBAGO COUNTY, ILLINOIS

Chemical of Potential Concern	Unit Risk	Units	Ac islment	Inhelation Cancer Slope Factor	Units	Weight of Evidence/ Cancer Guideline Description	Source	Date (1) (MM/DD/YY)
Acetone	N/A	N/A	N/A	N/A	N/A	D	IRIS	7/7/00
cis-1,2-Dichloroethene	N/A	N/A	N/A	N/A	N/A	D	IRIS	7/7/ <b>0</b> 0
Chloroform	2.30E-05	(ug/m3) <sup>-1</sup>	₹ <b>500</b>	8.10E-02	(mg/kg-day) <sup>-1</sup>	B2/hepatocellular carcinoma	IRIS	7/7/00
Benzene	2.2E-06 to 7.8E-06	(ug/m3) <sup>-1</sup>	3 <b>500</b>	7.7E-03 to 2.73E-02	(mg/kg-dey) <sup>-1</sup>	<b>Afleukemia</b>	IRIS	7/7/00
Trichloroethene	1.70E-06	(ug/m3) <sup>-1</sup>	3500	6.00E-03	(mg/kg-day) <sup>-t</sup>	N/A	NCEA/Region IX	11/29/99
Toluene	N/A	N/A	N/A	N/A	N/A	D	#RIS	7 <i>171</i> 00
Tetrachloroethene	5.70E-07	(ug/m3) <sup>-1</sup>	₹500	2.00E-03	(mg/kg-day) <sup>-1</sup>	liver	NCEA/Region IX	11/29/99
1,1,1-Trichloroethane	N/A	NA	N/A	N/A	N/A	D	IRIS	7/7/ <b>0</b> 0
1,1,2-Trichloro-1,2,2-trifluoromethane	N/A	N/A	NA	N/A	N/A	N/A B2/combined adenomas &	N/A	N/A
Methylene chloride	4.70E-07	(ug/m3) <sup>-1</sup>	3 OE+03	1.60E-03	(mg/kg-day) <sup>-1</sup>	carcinomas	IRIS	7/7/00
2-Butanone	N/A	N/A	NA	NA	N/A	. <b>D</b>	IRIS	7/7/00
m-&/or p-Xylene	N/A	N/A	N/A	N/A	NA	D	IRIS	7/7/00
Ethylbenzene	N/A	N/A	N/A	N/A	N/A	D	IRIS	7/7/00

(1) For IRIS values, provide the date IRIS was searched.

For HEAST values, provide the date of HEAST.

For NCEA values, provide the date of article provided by NCEA.

REGION IX = Region IX Preliminary remediation goals.

RPA = Relative Potency Approach (EPA, 1993s)

EPA Group:

- A Human carcinogen
- B1 Probable human carcinogen indicates that limited human data are available
- B2 Probable human carcinogen indicates sufficient evidence in animals and inadequate or no evidence in humans
- C Possible human carcinogen
- D Not classifiable as a human carcinogen
- E Evidence of noncarcinogenicity

Weight of Evidence:

Known/Likely

Cannot be Determined

**Not Likely** 

# TABLE 7.1.CT CALCUI ATION OF NON-CANCER HAZARDS CE ITRAL TENDENCY EXPOSURE EVERGREEN MANOF SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Medium: Exposure Medium: Current/Future Groundwater Groundwater

Exposure Point: Receptor Population: Tap water Resident

Receptor Age:

Adult

Exposure	Chemical	Medium	Medium	Route	Route	EPC	intake	Intake	Reference	Reference	Reference	Reference	Hazard
Route	of Potential	EPC	EPC	EPC	EPC	Selected	(Non-Cancer)	(Non-Cancer)	Dose (2)	Dose Units	Concentration	Concentration	Quotient
	Concern	Value	Units	Value	Units	for Hazard		Units				Units	
						Calculation (1)							
					_								
ngestion				400	_								
	Acetone	100	ug/L	100	ugL	M	1.918E-03	mg/kg-d	1.00E-01	mg/kg-d	-	-	1.918E-02
	Trichloroethene	6	ug/L	6	ug/L	. M	1.151E-04	mg/kg-d	6.00E-03	mg/kg-d	-	-	1.918E-02
	Tetrachioroethene	9	ug/L	9	ugiL	, M	1. <b>726E-04</b>	mg/kg-d	1.00E-02	mg/kg-d	-	-	1.726E-02
	Chloroform	0.9	ug/L	0.9	ug/L	M	1.726E-05	mg/kg-d	1.00E-02	mg/kg-d	_		1.726E-03
	Benzene	0.6	ug/L	0.6	ug/L	M	1.151E-05	mg/kg-d	3.00E-03	mg/kg-d	-	-	3.836E-03
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	9.589E-06	mg/kg-d	6.00E-02	mg/kg-d	-	-	1.598E-04
	(Total)		<u> </u>			ď.							6.1E-02
lermal absorption													
	Acetone	100	ug/L	100	ug/L	M	2.655E-06	mg/kg-d	1.00E-01	mg/kg-d	-	-	2.655E-05
	Trichioroethene	6	ug/L	6	ug/L	' М	4.471E-06	mg/kg-d	6.00E-03	mg/kg-d	-		7.452E-04
	Tetrachioroethene	9	ug/L	9	ug/L	M	2.012E-05	mg/kg-d	1.00E-02	mg/kg-d	-		2.012E-03
	Chloroform	0.9	ug/L	0.9	ug/L	, <b>M</b>	3.731E-07	mg/kg-d	1.00E-02	mg/kg-d	-	_	3.731E-05
	Benzene	0.6	ug/L	0.6	ugit.	ar M	5.868E-07	mg/kg-d	3.00E-03	mg/kg-d	-	_	1.956E-04
	Methylene chloride	0.5	ug/L	0.5	ugiL	M	1.048E-07	mg/kg-d	6.00E-02	mg/kg-d	_		1.747E-06
	(Total)				in the second	-							3.0E-03
halation													
	Acetone	100	ug/L	100	ug/L	M	1.027E-02	mg/kg-d	1.00E-01	mg/kg-d	1.0E-01	mg/kg-d	1.027E-01
	Trichloroethene	6	ug/L	6	ug/L	M	6.164E-04	mg/kg-d	6.00E-03	mg/kg-d	6.0E-03	mg/kg-d	1.027E-01
	Tetrachloroethene	9	ug/L	9	ug/L	M	9.247E-04	mg/kg-d	1.00E-02	mg/kg-d	1.1E-01	mg/kg-d	8.406E-03
	Chioroform	0.9	ug/L	0.9	ug/L	м	9.247E-05	mg/kg-d	1.00E-02	mg/kg-d	8.6E-05	mg/kg-d	1.075E+00
	Benzene	0.6	ug/L	0.6	ug/L	м	6.164E-05	mg/kg-d	3.00E-03	mg/kg-d	1.7E-03	mg/kg-d	3.626E-02
ļ	Methylene chloride	0.5	ug/L	0.5	ug/L	M	5.137E-05	mg/kg-d	6.00E-02	mg/kg-d	8.6E-01	mg/kg-d	5.973E-05
	(Total)	l		i									1.3E+00

Total Hazard Index Across All Exposure Routes/Pathways 1.4E+00

(1) Specify Medium-Specific (M) or Route-Specific (R) EPC selected for hazard calculation.

(2) Specify if subchronic.

N/A = Not Applicable

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#### TABLE 7.1.RME

#### CALCU: ATION OF NON-CANCER HAZARDS

#### REA JONABLE MAXIMUM EXPOSURE

EVERGREEN MANOF SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Current/Future
Medium: Groundwater
Exposure Medium: Groundwater
Exposure Point: Tap water
Receptor Population: Resident
Receptor Age: Adult

Tetrachi Chlorofo Benzene	loroethene chloroethene	Medium EPC Value	Medium EPC Units	Route EPC Value	Route EPC Units	EPC Selected for Hazard Calculation (1)	Intake (Non-Cancer)	Intake (Non-Cancer) Units	Reference Dose (2)	Reference Dose Units	Reference Concentration	Reference Concentration Units	Hazard Quotient
Acetone Trichlore Tetrachi Chlorofe Benzene Methylee	Concern one loroethene chloroethene	Value	Units			for Hazard	(Non-Cancer)		Dose (2)	Dose Units	Concentration		Quotient
Acetone Trichlon Tetrachi Chlorofo Benzene Methylei	one loroethene chloroethene			Value	Units	1		Units		1 1		1 1 1 1 1 1	
Acetone Trichlon Tetrachi Chlorofo Benzene Methylei	loroethene chloroethene	100				Calculation (1)				1 ?	1 1	Units	1
Acetone Trichlon Tetrachi Chlorofo Benzene Methylei	loroethene chloroethene	100											İ
Trichlon Tetrachi Chlorofo Benzeno Methylei	loroethene chloroethene	100									-		
Tetrachi Chlorofo Benzeno Methylei	chloroethene		ug/L	100	ug/L	M	2.740E-03	mg/kg-d	1.00E-01	mg/kg-d	] - [	-	2.740E-0
Chlorofo Benzeno Methyled		6	ug/L	6	ug/L	M ·	1.644E-04	mg/kg-d	6.00E-03	mg/kg-d	i - I	- 1	2.740E-0
Benzene Methylei		9	ug/L	9	ug/L	м	2.466E-04	mg/kg-d	1.00E-02	mg/kg-d	-	-	2.486E-02
Methylei	orom	0.9	ug/L	0.9	ug/L	м	2.466E-05	mg/kg-d	1.00E-02	mg/kg-d	-	- 1	2.466E-03
Permal absorption	ene	0.6	ug/L	0.6	ug/L	M	1.644E-05	mg/kg-d	3.00E-03	mg/kg-d	1	, ,	5.479E-03
· · · · · · · · · · · · · · · · · · ·	ylene chloride	0.5	ug/L	0.5	ug/L	M	1.370E-05	mg/kg-d	6.00E-02	mg/kg-d	] -		2.283E-04
· · · · · · · · · · · · · · · · · · ·	(Total)												8.6E-02
Acetone		,											
lucione.	ne	100	ug/L	100	ug/L	М	1.042E-05	mg/kg-d	1.00E-01	mg/kg-d	-	-	1.042E-04
Trichtore	oroethene	6	ug/L	6	ug/L	M	1.754E-05	mg/kg-d	6.00E-03	mg/kg-d	-	-	2.924E-03
Tetrachi	chloroethene	9	ug/L	9	ug/L	M	7.894E-05	mg/kg-d	1.00E-02	mg/kg-d	! - !	] - [	7.694E-03
Chlorofo	oform	0.9	ug/L	0.9	ug/L	М	1.464E-06	mg/kg-d	1.00E-02	mg/kg-d	-	-	1.464E-04
Benzene	ene	0.6	ug/L	0.6	ug/L	м	2.303E-06	mg/kg-d	3.00E-03	mg/kg-d	1 - 1	l - 1	7.675E-04
Methyler	lene chioride	0.5	ug/L	0.5	ug/L	М	4.112E-07	mg/kg-d	6.00E-02	mg/kg-d	-	-	6.853E-06
	(Total)												1.2E-02
halation			,										
Acetone	ne	100	ug/L	100	ug/L	м	1.027E-02	mg/kg-d	1.00E-01	mg/kg-d	1.0E-01	mg/kg-d	1.027E-01
Trichloro	proethene	6	ug/L	6	ug/L	м	6.164E-04	mg/kg-d	6.00E-03	mg/kg-d	6.0E-03	mg/kg-d	1.027E-01
Tetrachk	thloroethene	9	ug/L	9	ug/L	М	9.247E-04	mg/kg-d	1.00E-02	mg/kg-d	1.1E-01	mg/kg-d	8.406E-03
Chlorofo	oform	0.9	ug/L	0.9	ug/L	М	9.247E-05	mg/kg-d	1.00E-02	mg/kg-d	8.6E-05	mg/kg-d	1.075E+00
Benzene	ine	0.6	ug/L	0.6	ug/L	М	6.164E-05	mg/kg-d	3.00E-03	mg/kg-d	1.7E-03	mg/kg-d	3.626E-02
Methylen	lene chloride	0.5	ug/L	0.5	ug/L	M	5.137E-05	mg/kg-d	6.00E-02	mg/kg-d	8.6E-01	mg/kg-d	5.973E-05
	(Total)												

(1) Specify Medium-Specific (M) or Route-Specific (R) EPC selected for hazard calculation.

(2) Specify if subchronic.

N/A = Not Applicable

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### TABLE 7.2.CT

#### CALCU ATION OF NON-CANCER HAZARDS

#### CE ITRAL TENDENCY EXPOSURE

EVERGREEN MANOF SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Current/Future

Medium: Groundwater

Exposure Medium: Groundwater

Exposure Point: Tap water

Receptor Population: Resident

Receptor Age: Child

			1				1	1					
Exposure	Chemical	Medium	Medium	Route	Route	EPC	Intake	Intake	Reference	Reference	Reference	Reference	Hazard
Route	of Potential	EPC	EPC	EPC	EPC	Selected	(Non-Cancer)	(Non-Cancer)	Dose (2)	Dose Units	Concentration	Concentration	Quotier
	Concern	Value	Units	Value	Units	for Hazard		Units				Units	
		1		:		Calculation (1)							
	ļ. ————————————————————————————————————				-				ļ				
ngestion		100		100		м		mg/kg-d	1.00E-01	mg/kg-d			
	Acetone	6	ug/L	6	ug/L	, a	4.475E-03	mg/kg-d	6.00E-03	mg/kg-d	-	-	4.475E-
	Trichloroethene	_	ug/L		ug/L		2.685E-04			1 -	-	-	4.475E-
	Tetrachioroethene	9	ug/L	9	ug/L	× 4	4.027E-04	mg/kg-d	1.00E-02	mg/kg-d	-	-	4.027E-
	Chloroform	0.9	ug/L	0.9	ug/L		4.027E-05	mg/kg-d	1.00E-02	mg/kg-d			4.027E-
	Benzene	0.6 0.5	ug/L	0.6	ug/t.	M	2.685E-05	mg/kg-d	3.00E-03	mg/kg-d			8.950E-
	Methylene chloride	1	ug/L	0.5	ug/L	M	2.237E-05	mg/kg-d	6.00E-02	mg/kg-d	-	-	3.729E-
	(Total)								<u> </u>				1.4E-0
ermal absorption			i										
	Acetone	100	ug/L	100	ug/L	M	4.531E-06	mg/kg-d	1.00E-01	mg/kg-d	-	-	4.531E-
	Trichloroethene	6	ug/L	6	ug/L	м	7.631E-06	mg/kg-d	6.00E-03	mg/kg-d	-	-	1.272E-
	Tetrachloroethene	9	ug/L	9	ug/L	M	3.434E-05	mg/kg-d	1.00E-02	mg/kg-d	-	-	3.434E-
	Chioroform	0.9	ug/L	0.9	ug/L	M	6.367E-07	mg/kg-d	1.00E-02	mg/kg-d			6.367E-
	Benzene	0.6	ug/L	0.6	ug/L	M	1.002E-06	mg/kg-d	3.00E-03	mg/kg-d			3.338E-
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	1.788E-07	mg/kg-d	6.00E-02	mg/kg-d	-	-	2.981E-
	(Total)												5.2E-00
nhalation													
	Acetone	100	ug/L	100	ug/L	M	2.781E-02	mg/kg-d	1.00E-01	mg/kg-d	1.0E-01	mg/kg-d	2.781E-
	Trichloroethene	6	ug/L	6	ug/L	M	1.668E-03	mg/kg-d	6.00E-03	mg/kg-d	6.0E-03	mg/kg-đ	2.781E-
	Tetrachloroethene	9	ug/L	9	ugit	М	2.503E-03	mg/kg-d	1.00E-02	mg/kg-d	1.1E-01	mg/kg-d	2.275E-
	Chloroform	0.9	ug/L	0.9	ugl	M	2.503E-04	mg/kg-d	1.00E-02	mg/kg-d	8.6E-05	mg/kg-d	2.910E+
	Benzene	0.6	ug/L	0.6	ug/L	M	1.668E-04	mg/kg-d	3.00E-03	mg/kg-d	1.7E-03	mg/kg-d	9.815E-
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	1.390E-04	mg/kg-d	6.00E-02	mg/kg-d	8.6E-01	mg/kg-d	1.617E-
	(Total)				L								3.6E+0
	· · · · · · · · · · · · · · · · · · ·							Total Hazard	Index Acro	oss All Exp	sure Routes	/Pathways	3.7E+0

(1) Specify Medium-Specific (M) or Route-Specific (R) EPC selected for hazard calculation.

(2) Specify if subchronic.

N/A = Not Applicable

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#### TABLE 7.2.RME

#### CALCU: ATION OF NON-CANCER HAZARDS

#### REA ONABLE MAXIMUM EXPOSURE

EVERGREEN MANOF SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timetrame: Current/Future

Medium: Groundwater

Exposure Medium: Groundwater

Exposure Point: Tap water

Receptor Population: Resident

Receptor Age: Child

			1	i .				1					
Exposure Route	Chemical of Potential	Medium EPC	Medium EPC	Route EPC	Route EPC	EPC Selected	Intake (Non-Cancer)	Intake (Non-Cancer)	Reference Dose (2)	Reference Dose Units	Reference Concentration	Reference Concentration	Hazard
	Concern	Value	Units	Value	Units	for Hazard	,	Units	, ,			Units	
						Calculation (1)							
			1		1								
Ingestion							ì						
	Acetone	100	ug/L	100	ug/L	M	6.393E-03	mg/kg-d	1.00E-01	mg/kg-d	-	-	6.3 <b>93</b> E-0
	Trichloroethene	6	ug/L	6	ug/L	M	3.836E-04	mg/kg-d	6.00E-03	mg/kg-d	_	-	6.393E-0
	Tetrachioroethene	9	ug/L	9	ug/L	M	5.753E-04	mg/kg-d	1.00E-02	mg/kg-d	_	-	5.753E-0
	Chloroform	0.9	ug/L	0.9	ug/L	M	5.753E-05	mg/kg-d	1.00E-02	mg/kg-d	_		5.753E-0
	Benzene	0.6	ug/L	0.6	ug/L	M	3.836E-05	mg/kg-d	3.00E-03	mg/kg-d			1.279E-0
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	3.196E-05	mg/kg-d	6.00E-02	mg/kg-d	_ '	_	5.327E-0
	(Total)												2.0E-01
Dermal absorption													
	Acetone	100	ug/L	100	ug/L	М	1.804E-05	mg/kg-d	1.00E-01	mg/kg-d	l -	-	1.804E-0
	Trichloroethene	6	ug/L	6	ug/L	M	3.039E-05	mg/kg-d	6.00E-03	mg/kg-d	-		5.065E-0
	Tetrachloroethene	9	ug/L	9	ug/L	M	1.368E-04	mg/kg-d	1.00E-02	mg/kg-d	<b>-</b>	-	1.368E-0
	Chloroform	0.9	ug/L	0.9	ug/L	М	2.536E-06	mg/kg-d	1.00E-02	mg/kg-d		-	2.536E-0
	Benzene	0.6	ug/L	0.6	ug/L	M	3.969E-06	mg/kg-d	3.00E-03	mg/kg-d	-	-	1.330E-0
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	7.123E-07	mg/kg-d	6.00E-02	mg/kg-d	-		1.187E-0
	(Total)												2.1E-02
nhalation													
	Acetone	100	ug/L	100	ug/L	М	2.781E-02	mg/kg-d	1.00E-01	mg/kg-d	1.0E-01	mg/kg-d	2.781E-0
	Trichloroethene	6	ug/L	6	ug/L	M	1.668E-03	mg/kg-d	6.00E-03	mg/kg-d	6.0E-03	mg/kg-d	2.781E-0
	Tetrachioroethene	9	ug/L	9	ug/L	М	2.503E-03	mg/kg-d	1.00E-02	mg/kg-d	1.1E-01	mg/kg-d	2.275E-0
	Chloroform	0.9	ug/L	0.9	ug/L	м	2.503E-04	mg/kg-d	1.00E-02	mg/kg-d	8.6E-05	mg/kg-d	2.910E+0
	Benzene	0.6	ug/L	0.6	ug/L	м	1.668E-04	mg/kg-d	3.00E-03	mg/kg-d	1.7E-03	mg/kg-d	9.815E-0
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	1.390E-04	mg/kg-d	6.00E-02	rng/kg-d	8.6E-01	mg/kg-d	1.617E-0
	(Total)												3.6E+00
			· · · · · · · · · · · · · · · · · · ·		_			Total Hazard	Index Acro	ss All Exp	sure Routes	/Pathways	3.8E+0

(1) Specify Medium-Specific (M) or Route-Specific (R) EPC selected for hazard calculation.

(2) Specify if subchronic.

N/A = Not Applicable

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#### TABLE 7.3.CT CALCUI ATION OF NON-CANCER HAZARDS CE ITRAL TENDENCY EXPOSURE

EVERGREEN MANOF SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe:

Future

Medium:

Groundwater

Exposure Medium:

Groundwater

Exposure Point:

Tap water

Receptor Population:

Commercial/Industrial

Receptor Age:

Adult

Exposure Route	Chemical of Potential Concern	Medium EPC Value	Medium EPC Units	Route EPC Value	Route EPC Units	EPC Selected for Hazard Calculation (1)	Intake (Non-Cancer)	Intake (Non-Cancer) Units	Reference Dose (2)	Reference Dose Units	Reference Concentration	Reference Concentration Units	Hazard Quotient
Ingestion		1						1					
	Acetone	100	ug/L	100	ug/L	м	6.849E-04	mg/kg-d	1.00E-01	mg/kg-d	_	-	6.849E-03
	Trichloroethene	6	ug/L	6	ug/L	М	4.110E-05	mg/kg-d	6.00E-03	mg/kg-d	_	-	6.849E-03
	Tetrachloroethene	9	ug/L	9	ug/L	M	6.164E-05	mg/kg-d	1.00E-02	mg/kg-d	- !		6.164E-03
	Chioroform	0.9	ug/L	0.9	ug/L	M	6.164E-06	mg/kg-d	1.00E-02	mg/kg-d	-	_	6.164E-04
1	Benzene	0.6	ug∕L	0.6	ug/L	M	4.110E-06	mg/kg-d	3.00E-03	mg/kg-d	_	-	1.370E-03
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	3.425E-06	mg/kg-d	6.00E-02	mg/kg-d	[ - ]	-	5.708E-05
	(Total)												2.2E-02
Dermal absorption													
	Acetone	100	ug/L	100	ug/L	M	1.896E-06	mg/kg-d	1.00E-01	mg/kg-d	_	-	1.896E-05
	Trichloroethene	6	ug/L	6	ug/L	M	3.194E-06	mg/kg-d	6.00E-03	mg/kg-d	_	_	5.323E-04
	Tetrachloroethene	9	ug/L	9	ug/L	M	1.437E-05	mg/kg-d	1.00E-02	mg/kg-d	-	_	1.437E-03
	Chloroform	0.9	ug/L	0.9	ug/L	М	2.665E-07	mg/kg-d	1.00E-02	mg/kg-d	-	-	2.665E-05
	Benzene	0.6	ug/L	0.6	ug/L	M	4.192E-07	mg/kg-d	3.00E-03	mg/kg-d	-	_	1.397E-04
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	7.485E-08	mg/kg-d	6.00E-02	mg/kg-d	-		1.248E-06
	(Total)												2.2E-03
Inhalation													
	Acetone	100	ug/L	100	ug/L	M	7.339E-03	mg/kg-d	1.00E-01	mg/kg-d	1.0E-01	mg/kg-d	7.339E-02
	Trichloroethene	6	ug/L	6	ug/L	M	4.403E-04	mg/kg-d	6.00E-03	mg/kg-d	6.0E-03	mg/kg-d	7.339E-02
	Tetrachioroethene	9	ug/L	9	ugit	M	6.605E-04	mg/kg-d	1.00E-02	mg/kg-d	1.1E-01	mg/kg-d	6.004E-03
	Chloroform	0.9	ug/L	0.9	ug/L	M	6.605E-05	mg/kg-d	1.00E-02	mg/kg-d	8.6E-05	mg/kg-d	7.680E-01
	Benzene	0.6	ug/L	0.6	ug/L	М	4.403E-05	mg/kg-d	3.00E-03	mg/kg-d	1.7E-03	mg/kg-d	2.590E-02
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	3.669E-05	mg/kg-d	6.00E-02	mg/kg-d	8.6E-01	mg/kg-d	4.267E-05
	(Total)									L			9.5E-01
								Total Hazard	Index Acre	oss All Expe	sure Routes	/Pathways	9.7E-01

(1) Specify Medium-Specific (M) or Route-Specific (R) EPC selected for hazard calculation.

(2) Specify if subchronic.

N/A = Not Applicable

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#### TABLE 7.3.RME

#### CALCU ATION OF NON-CANCER HAZARDS

#### REA JONABLE MAXIMUM EXPOSURE

EVERGREEN MANO: SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Future
Medium: Groundwater
Exposure Medium: Groundwater

Exposure Point: Tap water

Receptor Population: Commercial/Industrial Receptor Age: Adult

		7	1		-	1		1		· · · · · · · · · · · · · · · · · · ·			
Exposure Route	Chemical of Potential	Medium EPC	Medium EPC	Route EPC	Route EPC	EPC Selected	Intake (Non-Cancer)	Intake (Non-Cancer)	Reference Dose (2)	Reference Dose Units	Reference Concentration	Reference Concentration	Hazard Quotier
	Concern	Value	Units	Value	Units	for Hazard		Units	``			Units	
	J	1				Calculation (1)			1				
		j				.,,							
Ingestion					_					Ì			
	Acetone	100	ug/L	100	ug/L	M	9.785E-04	mg/kg-d	1.00E-01	mg/kg-d	_	_	9.785E-0
	Trichloroethene	6	ug/L	6	ug/L	M	5.871E-05	mg/kg-d	6.00E-03	mg/kg-d	_	-	9.785E-0
	Tetrachloroethene	9	ug/L	9	ug/L	) M ]	8.806E-05	mg/kg-d	1.00E-02	mg/kg-d	ì -	] -	8.806E-0
	Chloroform	0.9	ug/L	0.9	ug/L	M	8.806E-06	mg/kg-d	1.00E-02	mg/kg-d	-	-	8.806E-0
	Benzene	0.6	ug/L	0.6	ug/L	M	5.871E-06	mg/kg-d	3.00E-03	mg/kg-d	-	-	1.957E-0
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	4.892E-06	mg/kg-d	6.00E-02	mg/kg-d	_	- :	8.154E-0
	(Total)					[							3.1E-02
Dermal absorption													-
	Acetone	100	ug/L	100	ug/L	м	7.440E-06	mg/kg-d	1.00E-01	mg/kg-d	-	-	7.440E-0
	Trichloroethene	6	ug/L	6	ug/L	М	1.253E-05	mg/kg-d	6.00E-03	mg/kg-d	_	-	2.088E-0
	Tetrachloroethene	9	ug/L	9	ug/L	м	5.639E-05	mg/kg-d	1. <b>00E-0</b> 2	mg/kg-d	-	-	5.639E-0
	Chioroform	0.9	ug/L	0.9	ug/L	M	1.046E-06	mg/kg-d	1.00E-02	mg/kg-d	-	l -	1.046E-0
	Benzene	0.6	ug/L	0.6	ug/L	м	1.645E-06	mg/kg-d	3.00E-03	mg/kg-d	-	-	5.482E-0
	Methylene chloride	0.5	ug/L	0.5	ug/L	М	2.937E-07	mg/kg-d	6.00E-02	mg/kg-d	-	-	4.895E-0
	(Total)												8.5E-03
nhalation													
	Acetone	100	ug/L	100	ug/L	M	7.339E-03	mg/kg-d	1.00E-01	mg/kg-d	1.0E-01	mg/kg-d	7.339E-0
	Trichloroethene	6	ug/L	6	ug/L	М	4.403E-04	mg/kg-d	6.00E-03	mg/kg-d	6.0E-03	mg/kg-d	7.339E-0.
	Tetrachioroethene	9	ug/L	9	ug/L	M	6.605E-04	mg/kg-d	1.00E-02	mg/kg-d	1.1E-01	mg/kg-d	6.004E-0
	Chloroform	0.9	ug/L	0.9	ug/L	M	6.605E-05	mg/kg-d	1.00E-02	mg/kg-d	8.6E-05	mg/kg-d	7.680E-0
	Benzene	0.6	ug/L	0.6	ug/L	M	4.403E-05	mg/kg-d	3.00E-03	mg/kg-d	1.7E-03	mg/kg-d	2.590E-0
	Methylene chloride	0.5	ug/L	0.5	ng/L	M	3.889E-05	mg/kg-d	6.00E-02	mg/kg-d	8.6E-01	mg/kg-d	4.267E-0
	(Total)			لحجيب									9.5E-01
								Total Hazard	Index Acro	oss All Expo	sure Routes	/Pathways	9.9E-01

(1) Specify Medium-Specific (M) or Route-Specific (R) EPC selected for hazard calculation.

(2) Specify if subchronic.

N/A = Not Applicable

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#### TABLE 6.1.CT

#### CALCULATION OF CANCER RISKS

#### CENTRAL TENDENCY EXPOSURE

EVERGREEN MANO: SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Current/Future

Medium: Groundwater

Exposure Medium: Groundwater

Exposure Point: Tap Water

Receptor Population: Resident

Receptor Age: Adult

Exposure	Chemical	Medium	Medium	Route	Route	EPC Selected	Intake	intake	Cancer Slope	Cancer Stope	Cancer
Route	of Potential	EPC	EPC	EPC	EPC	for Risk	(Cancer)	(Cancer)	Factor	Factor Units	Risk
	Concern	Value	Units	Value	. Units	Calculation (1)		Units			
ngestion				-						<u> </u>	
	Acetone	100	ug/L	100	ugit.	М	1.918E-04	mg/kg-d	NC	mg/kg-d <sup>-1</sup>	-
	Trichloroethene	6	ug/L	6	ug/L	M	1.151E-05	mg/kg-d	1.10E-02	mg/kg-d <sup>-1</sup>	1.27E-07
	Tetrachloroethene	9	ug/L	9	ugit	M	1.726E-05	mg/kg-d	5.20E-02	mg/kg-d <sup>-1</sup>	8.98E-07
	Chloroform	0.9	ug/L	0.9	ugit.	М	1.726E-06	mg/kg-d	6.10E-03	mg/kg-d <sup>-1</sup>	1.05E-08
	Benzene	0.6	ug/L	0.6	ugf.	M	1.151E-06	mg/kg-d	5.50E-02	mg/kg-d <sup>-1</sup>	6.33E-06
	Methylene chloride	0.5	ug/L	0.5	ugit.	M	9.589E-07	mg/kg-d	7.50E-03	mg/kg-d <sup>-1</sup>	7.19E-09
	(Total)										1.11E-06
ermal											
	Acetone	100	ug∕L	100	ugil,	M	2.655E-07	mg/kg-d	NC	mg/kg-d <sup>-1</sup>	-
	Trichioroethene	6	ug/L	6	ug/L,	M	4.471E-07	mg/kg-d	1.10E-02	mg/kg-d <sup>-1</sup>	4.92E-09
	Tetrachloroethene	9	ug/L	9	ug/L	M	2.012E-06	mg/kg-d	5.20E-02	mg/kg-d <sup>-1</sup>	1.05E-07
	Chloroform	0.9	ug/L	0.9	ugf.	) M	3.731E-08	mg/kg-d	6.10E-03	mg/kg-d <sup>-1</sup>	2.28E-10
	Benzene	0.6	ug/L	0.6	ugli.	M	5.868E-08	mg/kg-d	5.50E-02	mg/kg-d <sup>-1</sup>	3.23E-09
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	1.048E-08	mg/kg-d	7.50E-03	mg/kg-d <sup>-1</sup>	7.86E-11
	(Total)			-							1.13E-07
halation										_	
	Acetone	100	ug/L	100	ug/L	M	1.027E-03	mg/kg-d	NC	mg/kg-d <sup>-1</sup>	
	Trichloroethene	6	ug/L	6	mpl.	M	6.164E-05	mg/kg-d	6.00E-03	mg/kg-d <sup>-1</sup>	3.70E-07
	Tetrachloroethene	9	ug/L	•	angle.	M	9.247E-05	mg/kg-d	2.00E-03	mg/kg-d <sup>-1</sup>	1.85E-07
i	Chloroform	0.9	ug/L	0.9	upt.	M	9.247E-06	mg/kg-d	8.10E-02	mg/kg-d <sup>-1</sup>	7.49E-07
	Benzene	0.6	ug/L	0.6	ugl.	M	6.164E-06	mg/kg-d	2.73E-02	mg/kg-d <sup>-1</sup>	1.66E-07
	Methylene chloride	0.5	ug/L	0.5	ugt	М	5.137E-06	mg/kg-d	1.60E-03	mg/kg-d <sup>-1</sup>	8.22E-09
	(Total)	{							j i		1.48E-08

(1) Specify Medium-Specific (M) or Route-Specific (R) EPC selected for risk calculation.

N/A = Not Applicable

NC = Not carcinogenic.

#### TABLE 8.1.RME

#### C/ CULATION OF CANCER RISKS

#### RE/ SONABLE MAXIMUM EXPOSURE

EVERGREEN MANO : SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Current/Future
Medium: Groundwater
Exposure Medium: Groundwater
Exposure Point: Tap Water

Receptor Population: Resident Receptor Age: Adult

Exposure	Chemical	Medium	Medium	Route	Route	EPC Selected	Intake	Intake	Cancer Slope	Cancer Slope	Cancer
Route	of Potential	EPC	EPC	EPC	EPC	for Risk	(Cancer)	(Cancer)	Factor	Factor Units	Risk
	Concern	Value	Units	Value	Units	Calculation (1)		Units			
ngestion				-	<u> </u>			<u> </u>			
	Acelone	100	ug/L	100	ug/L	M	9.393E-04	mg/kg-d	NC	mg/kg-d <sup>-1</sup>	-
	Trichloroethene	6	ug/L	6	ug/L	M	5.636E-05	mg/kg-d	1.10E-02	mg/kg-d <sup>-1</sup>	6.20E-07
	Tetrachloroethene	9	ug/L	9	ug/L	M	8.454E-05	mg/kg-d	5.20E-02	mg/kg-d <sup>-1</sup>	4.40E-08
	Chloroform	0.9	ug/L	0.9	ug/L	) M	8.454E-06	mg/kg-d	6.10E-03	mg/kg-d <sup>-1</sup>	5.16E-08
	Benzene	0.6	ug/L	0.6	ug/L	M	5.636E-06	mg/kg-d	5.50E-02	mg/kg-d <sup>-1</sup>	3.10E-07
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	4.897E-06	mg/kg-d	7.50E-03	mg/kg-d <sup>-1</sup>	3.52E-08
	(Total)										5.41E-06
ermal				-							
	Acetone	100	ug/L	100	ug/L	M	3.571E-06	mg/kg-d	NC	mg/kg-d <sup>-1</sup>	
	Trichloroethene	6	ug/L	6	ug/L	М	6.015E-06	mg/kg-d	1.10E-02	mg/kg-d <sup>-1</sup>	6.62E-06
	Tetrachloroethene	9	ug/L	8	ug/L	M	2.707E-05	mg/kg-d	5.20E-02	mg/kg-d <sup>-1</sup>	1.41E-06
	Chloroform	0.9	ug/L	0.9	ug/L	M	5.019E-07	mg/kg-d	6.10E-03	mg/kg-d <sup>-1</sup>	3.06E-09
	Benzene	0.6	ug/L	0.6	ug/L	M	7.894E-07	mg/kg-d	5.50E-02	mg/kg-d <sup>-1</sup>	4.34E-08
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	1.410E-07	mg/kg-d	7.50E-03	mg/kg-d <sup>-1</sup>	1.06E-09
	(Total)										1.52E-06
halation											
	Acetone	100	ug/L	100	ug/L	, M	3.523E-03	mg/kg-d	NC	mg/kg-d <sup>-1</sup>	-
	Trichloroethene	6	ug/L	6	ug/L	M	2.114E-04	mg/kg-d	6.00E-03	mg/kg-d <sup>-1</sup>	1.27E-06
	Tetrachloroethene	9	ug/L	9	ug/L	M	3.170E-04	mg/kg-d	2.00E-03	mg/kg-d <sup>-1</sup>	6.34E-07
	Chloroform	0.9	ug/L	0.9	ug/L	М	3.170E-05	mg/kg-d	8.10E-02	mg/kg-d <sup>-1</sup>	2.57E-08
	Benzene	0.6	ug/L	0.6	ug/L	M	2.114E-05	mg/kg-d	2.73E-02	mg/kg-d <sup>-1</sup>	5.77E-07
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	1.761E-05	mg/kg-d	1.60E-03	mg/kg-d <sup>-1</sup>	2.82E-08
i	(Total)		1						1	ſ	5.08E-06

(1) Specify Medium-Specific (M) or Route-Specific (R) EPC selected for risk calculation.

N/A = Not Applicable

NC = Not carcinogenic.

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#### TABLE 8.2.CT

#### CA CULATION OF CANCER RISKS

#### CE ITRAL TENDENCY EXPOSURE

EVERGREEN MANO: SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Current/Future
Medium: Groundwater
Exposure Medium: Groundwater
Exposure Point: Tap Water
Receptor Population: Resident
Receptor Age: Child

Expasure Route	Chemical of Potential Concern	Medium EPC Value	Medium EPC Units	Route EPC Value	Route EPC Units	EPC Selected for Risk Calculation (1)	intake (Cancer)	intake (Cancer) Units	Cancer Slope Factor	Cancer Slope Factor Units	Cancer Risk
Ingestion											
	Acetone	100	ug/L	100	ug/L	M	1.279E-04	mg/kg-d	NC	mg/kg-d <sup>-1</sup>	-
	Trichloroethene	6	ug/L	6	ug/i.	M	7.871E-08	mg/kg-d	1.10E-02	mg/kg-d <sup>-1</sup>	8.44E-08
	Tetrachloroethene	9	ug/L	9	ug/L	M	1.151E-05	mg/kg-d	5.20E-02	mg/kg-d <sup>-1</sup>	5.98E-07
	Chloroform	0.9	ug/L	0.9	ugl.	М	1.151E-06	mg/kg-d	6.10E-03	mg/kg-d <sup>-1</sup>	7.02E-09
	Benzene	0.6	ug/L	0.6	ugl	м	7.871E-07	mg/kg-d	5.50E-02	mg/kg-d <sup>-1</sup>	4.22E-08
	Methylene chloride	0.5	ug/L	0.5	ugA.	м	6.393E-07	mg/kg-d	7.50E-03	mg/kg-d <sup>-1</sup>	4.79E-09
	(Total)	<u> </u>									7.37E-07
Dermal											
	Acetone	100	.ug/L	100	ug/L	M	1.294E-07	mg/kg-d	NC .	mg/kg-d <sup>-1</sup>	-
	Trichloroethene	6	ug/L	6	ugiL	M	2.180E-07	mg/kg-d	1.10E-02	mg/kg-d <sup>-1</sup>	2.40E-09
	Tetrachloroethene	9	ug/L	9	ugt.	M	9.811E-07	mg/kg-d	5.20E-02	mg/kg-d <sup>-1</sup>	5.10E-08
	Chloroform	0.9	ug/L	0.9	agt.	M	1.819E-08	mg/kg-d	6.10E-03	mg/kg-d <sup>-1</sup>	1.11E-10
	Benzene	0.6	ug/L	0.6	ugit.	M	2.861E-06	mg/kg-d	5.50E-02	mg/kg-d <sup>-1</sup>	1.57E-09
	Methylene chloride	0.5	ug/L	0.5	ugiL	M	5.110E-09	mg/kg-d	7.50E-03	mg/kg-d <sup>-1</sup>	3.83E-11
	(Total)				ı						5.51E-08
nhalation											
	Acetone	100	ug/L	100	ugl	M	7.945E-04	mg/kg-d	NC NC	mg/kg-d <sup>-1</sup>	-
	Trichloroethene	6	ug/L	6	<b>apt</b>	M	4.767E-05	mg/kg-d	6.00E-03	mg/kg-d <sup>-1</sup>	2.86E-07
	Tetrachloroethene	9	ug/L	9	egt.	M	7.151E-05	mg/kg-d	2.00E-03	mg/kg-d <sup>-1</sup>	1.43E-07
	Chloroform	0.9	ug/L	0.9	ugiL	М	7.151E-06	mg/kg-d	8.10E-02	mg/kg-d <sup>-1</sup>	5.79E-07
	Benzene	0.6	ug/L	0.6	ug/L	М	4.767E-06	mg/kg-d	2.73E-02	mg/kg-d <sup>-1</sup>	1.30E-07
	Methylene chloride	0.5	ug/L	. 0.5	ug/L	М	3.973E-06	mg/kg-d	1.60E-03	mg/kg-d <sup>-1</sup>	6.36E-09
	(Total)										1.14E-08
						Total H	azard Index	Across All Ex	coosure Route	s/Pathways	1.94E-06

(1) Specify Medium-Specific (M) or Route-Specific (R) EPC selected for risk calculation.

N/A = Not Applicable

NC = Not carcinogenic.

#### TABLE 8.2.RME

#### CA CULATION OF CANCER RISKS

#### REA JONABLE MAXIMUM EXPOSURE

#### EVERGREEN MANO: SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Current/Future ledium: Groundwater Exposure Medium: Groundwater Exposure Point: Tap Water Receptor Population: Resident Receptor Age: Child

					•						
Exposure	Chemical	Medium	Medium	Route	Route	EPC Selected	Intake	Intake	Cancer Slope	Cancer Slope	Cancer
Route	of Potential	EPC	EPC	EPC	EPC	for Risk	(Cancer)	(Cancer)	Factor	Factor Units	Risk
	Concern	Value	Units	Value	Units	Calculation (1)		Units			
ngestion							· · · · · · · · · · · · · · · · · · ·	<del> </del>		<del>,</del>	
	Acetone	100	ug/L	100	ug/L	м	5.479E-04	mg/kg-d	NC	mg/kg-d <sup>-1</sup>	-
	Trichloroethene	6	ug/L	6	ug/L	M	3.288E-05	mg/kg-d	1.10E-02	mg/kg-d <sup>-1</sup>	3.62E-0
	Tetrachloroethene	9	ug/L	9	ug/L	M	4.932E-05	mg/kg-d	5.20E-02	mg/kg-d <sup>-1</sup>	2.56E-0
	Chloroform	0.9	ug/L	0.9	ug/L	M	4.932E-06	mg/kg-d	6.10E-03	mg/kg-d <sup>-1</sup>	3.01E-0
	Benzene	0.6	ug/L	0.6	ug/L	M	3.288E-06	mg/kg-d	5.50E-02	mg/kg-d <sup>-1</sup>	1.81E-0
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	2.740E-08	mg/kg-d	7.50E-03	mg/kg-d <sup>-1</sup>	2.05E-0
	(Total	)			L						3.16E-0
Permal				-							
	Acetone	100	ug/L	100	ug/L	M	1.547E-08	mg/kg-d	NC NC	mg/kg-d <sup>-1</sup>	-
	Trichloroethene	6	ug/L	6	ug/L	M	2.605E-08	mg/kg-d	1.10E-02	mg/kg-d <sup>-1</sup>	2.87E-0
	Tetrachloroethene	9	ug/L	9	ug/L	M	1.172E-05	mg/kg-d	5.20E-02	mg/kg-d <sup>-1</sup>	6.10E-0
	Chloroform	0.9	ug/L	0.9	ug/L	M	2.173E-07	mg/kg-d	6.10E-03	mg/kg-d <sup>-1</sup>	1.33E-0
	Benzene	0.6	ug/L	0.6	ug/L	M	3.419E-07	mg/kg-d	5.50E-02	mg/kg-d <sup>-1</sup>	1.88E-0
	Methylene chloride	0.5	ug/L	0.5	ug/L	M	6.105E-08	mg/kg-d	7.50E-03	mg/kg-d <sup>-1</sup>	4.58E-1
	(Total	)									6.59E-0
halation											
	Acetone	100	ug/L	100	ug/L	M	2.384E-03	mg/kg-d	NC	mg/kg-d <sup>-1</sup>	-
	Trichloroethene	6	ug/L	8	ug/L	M	1.430E-04	mg/kg-d	6.00E-03	mg/kg-d <sup>-1</sup>	8.58E-0
	Tetrachioroethene	9	ug/l.	9	ug/L	M	2.145E-04	mg/kg-d	2.00E-03	mg/kg-d <sup>-1</sup>	4.29E-0
	Chloroform	0.9	ug/L	0.9	ug/L	M	2.145E-05	mg/kg-d	8.10E-02	mg/kg-d <sup>-1</sup>	1.74E-0
	Benzene	0.6	ug/L	0.6	ug/L	M	1.430E-05	mg/kg-d	2.73E-02	mg/kg-d <sup>-1</sup>	3.90E-0
	Methylene chloride	0.5	ug/t_	0.5	ug/L	M	1.192E-05	mg/kg-d	1.60E-03	mg/kg-d <sup>-1</sup>	1.91E-0
	(Total										3.43E-0
			-			Total H	lazard Index	Across All E	cposure Route	s/Pathways	7.25E-0

(1) Specify Medium-Specific (M) or Route-Specific (R) EPC selected for risk calculation.

N/A = Not Applicable NC = Not carcinogenic.

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#### TABLE 8.3.CT

#### CA CULATION OF CANCER RISKS

#### CE ITRAL TENDENCY EXPOSURE

#### EVERGREEN MANO: SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe:

Future

Medium:

Groundwater

Exposure Medium:

Groundwater Tap Water

Exposure Point: Receptor Population:

Commercial/Industrial

Receptor Age:

Adult

Exposure	Chemical	Medium	Medium	Route	Route	EPC Selected	Intake	intake	Cancer Slope	Cancer Slope	Cancer
Route	of Potential	EPC	EPC	EPC	EPC	for Risk	(Cancer)	(Cancer)	Factor	Factor Units	Risk
	Concern	Value	Units	Value	Units	Calculation (1)		Units			
Ingestion		· · · · · · · · · · · · · · · · · · ·					·				
	Acetone	100	ug/L	100	ug/L	M	8.806E-05	mg/kg-d	NC NC	mg/kg-d <sup>-1</sup>	-
	Trichioroethene	6	ug/L	6	ug/L	M	5.284E-06	mg/kg-d	1.10E-02	mg/kg-d <sup>-1</sup>	5.81E-08
	Tetrachioroethene	9	ug/L	9	ug/L	M	7.926E-06	mg/kg-d	5.20E-02	mg/kg-d <sup>-1</sup>	4.12E-07
	Chloroform	0.9	ug/L	0.9	ug/L	M	7.926E-07	mg/kg-d	6.10E-03	mg/kg-d <sup>-1</sup>	4.83E-09
	Benzene	0.6	ug/L	0.6	ug/L	M	5.284E-07	mg/kg-d	5.50E-02	mg/kg-d <sup>-1</sup>	2.91E-08
	Methylene chloride	0.5	ug/L	0.5	upl	M	4.403E-07	mg/kg-d	7.50E-03	mg/kg-d <sup>-1</sup>	3.30E-09
	(Total)						L		<u></u>		5.07E-07
Dermal											
	Acetone	100	ug/L	100	ug/L	M	2.438E-07	mg/kg-d	NC	mg/kg-d <sup>-1</sup>	-
	Trichloroethene	6	ug/L	6	ug/L	M	4.106E-07	mg/kg-d	1.10E-02	mg/kg-d <sup>-1</sup>	4.52E-09
	Tetrachioroethene	9	ug/L	9	ug/L	M	1.848E-06	mg/kg-d	5.20E-02	mg/kg-d <sup>-1</sup>	9.61E-08
	Chloroform	0.9	ug/L	0.9	ugit	M	3.426E-08	mg/kg-d	6.10E-03	mg/kg-d <sup>-1</sup>	2.09E-10
	Benzene	0.6	υ <b>g/L</b>	0.6	ugiL	M	5.389E-08	mg/kg-d	5.50E-02	mg/kg-d <sup>-1</sup>	2.96E-09
	Methylene chloride	0.5	ug/L	0.5	ug/L	м	9.624E-09	mg/kg-d	7.50E-03	mg/kg-d <sup>-1</sup>	7.22E-11
	(Total)			-							1.04E-07
Inhalation		400	_	400	7584 -						
	Acetone	100	ug/L	100	ugL	M	9.435E-04	mg/kg-d	NC	mg/kg-d <sup>-1</sup>	
	Trichloroethene	6	ug/L	6	ugL	M	5.661E-05	mg/kg-d	6.00E-03	mg/kg-d <sup>-1</sup>	3.40E-07 1.70E-07
	Tetrachioroethene	9	ug/L	9	ugit	M	8.492E-05	mg/kg-d	2.00E-03	mg/kg-d <sup>-1</sup>	1.70E-07 6.88E-07
	Chloroform	0.9	ug/L	0.9	ug/L	M M	8.492E-06	mg/kg-d	8.10E-02	mg/kg-d <sup>-1</sup>	6.88E-07 1.55E-07
	Benzene	0.6 0.5	ug/L	0.6 0.5	ugl	🖁	5.661E-06	mg/kg-d	2.73E-02	mg/kg-d <sup>-1</sup> mg/kg-d <sup>-1</sup>	1.55E-07 7.55E-09
	Methylene chloride	***	ug/L	U.J	ug/L	[ <b>"</b>	4.718E-06	mg/kg-d	1.60E-03	indaya-a	
	(Total)			-	<u> L</u>		لجببا				1.36E-06
						Total H	azard Index	Across All Ex	posure Route	es/Pathways	1.97E-06

(1) Specify Medium-Specific (M) or Route-Specific (R) EPC selected for risk calculation.

N/A = Not Applicable

NC = Not carcinogenic.

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#### TABLE 8.3.RME

#### C/ .CULATION OF CANCER RISKS

RE/ SONABLE MAXIMUM EXPOSURE

EVERGREEN MANO SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe:

Future

Medium:

Groundwater

Exposure Medium:

Groundwater

Exposure Point:

Tap Water

Receptor Population:

Commercial/Industrial

Receptor Age:

Adult

	<u> </u>		1				Y				
Exposure	Chemical	Medium	Medium	Route	Route	EPC Selected	intake	Intake	Cancer Slope	Cancer Slope	Cancer
Route	of Potential	EPC	EPC	EPC	EPC	for Risk	(Cancer)	(Cancer)	Factor	Factor Units	Risk
	Concern	Value	Units	Value	Units	Calculation (1)		Units			
Ingestion					1						
	Acetone	100	ug/L	100	ug/L	М	3.495E-04	mg/kg-d	NC NC	mg/kg-d <sup>-1</sup>	-
	Trichloroethene	6	ug/L	6	ug/L	м	2.097E-05	mg/kg-d	1.10E-02	mg/kg-d <sup>-1</sup>	2.31E-07
	Tetrachioroethene	9	ug/L	9	ug/L	М	3.145E-05	mg/kg-d	5.20E-02	mg/kg-d <sup>-1</sup>	1.64E-06
	Chloroform	0.9	ug/L	0.9	ug/L	M	3.145E-06	mg/kg-d	6.10E-03	mg/kg-d <sup>-1</sup>	1.92E-08
	Benzene	0.6	ug/L	0.6	ug/L	м	2.097E-06	mg/kg-d	5.50E-02	mg/kg-d <sup>-1</sup>	1.15E-07
	Methylene chloride	0.5	ug/L	0.5	ug/L	М	1.747E-06	mg/kg-d	7.50E-03	mg/kg-d <sup>-1</sup>	1.31E-08
<u> </u>	(Total)				1						2.01E-06
Dermal									i		
İ	Acetone	100	ug/L	100	ug/L	M	2.657E-06	mg/kg-d	NC NC	mg/kg-d <sup>-1</sup>	-
ľ	Trichloroethene	6	ug∕L	6	ug/L	M	4.475E-06	mg/kg-d	1.10E-02	mg/kg-d <sup>-1</sup>	4.92E-06
ļ	Tetrachioroethene	9	ug∕L	9	ug/L	M	2.014E-05	mg/kg-d	5.20E-02	mg/kg-d <sup>-1</sup>	1.05E-06
ļ.	Chloroform	0.9	ug/L	0.9	ug/L	M	3.734E-07	mg/kg-d	6.10E-03	mg/kg-d <sup>-1</sup>	2.28E-09
ļ	Benzene	0.6	ug/L	0.6	ug/L	M	5.874E-07	mg/kg-d	5.50E-02	mg/kg-d <sup>-1</sup>	3.23E-08
· 1	Methylene chloride	0.5	ug/L	0.5	ug/L	M	1.049E-07	mg/kg-d	7.50E-03	mg/kg-d <sup>-1</sup>	7.87E-10
	(Total)									<del></del>	1.13E-08
Inhelation											
	Acetone	100	ug/L	100	ug/L	M	2.621E-03	mg/kg-d	NC NC	mg/kg-d <sup>-1</sup>	
<b>i</b>	Trichloroethene	6	ug/L	6	ug/L	M	1.573E-04	mg/kg-d	6.00E-03	mg/kg-d <sup>-1</sup>	9.44E-07
ľ	Tetrachioroethene	9	ug/L	9	ug/L	M	2.359E-04	mg/kg-d	2.00E-03	mg/kg-d <sup>-1</sup>	4.72E-07
1	Chloroform	0.9	ug/L	0.9	ug/L	M	2.359E-05	mg/kg-d	8.10E-02	mg/kg-d <sup>-1</sup>	1.91E-06
i í	Benzene	0.6	ug/L	0.6	ug/L	М	1.573E-05	mg/kg-d	2.73E-02	mg/kg-d <sup>-1</sup>	4.29E-07
	Methylene chloride	0.5	ug/L	0.5	ug/L	м	1.310E-05	mg/kg-d	1.60E-03	mg/kg-d <sup>-1</sup>	2.10E-08
	(Total)										3.78E-06
		·				Total H	azard Index	Across All Ex	posure Route	s/Pathways	6.92E-08

(1) Specify Medium-Specific (M) or Route-Specific (R) EPC selected for risk calculation.

N/A = Not Applicable

NC = Not carcinogenic.

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#### TABLE 9.1.CT

SUMMARY OF FECEPTOR RISKS AND HAZARDS FOR COPCs

C NTRAL TENDENCY EXPOSURE

EVERGREEN MANO : SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timefram Current/Future Receptor Populatio Resident Receptor Age: Adult

Medium	Exposure Medium	Exposure Point	Chemical	<del></del>	C. cir	nogenic Risk		Chemical		Non-Carcin	ogenic Hazard	Quotient	
			İ	Ingestion	Inhalati n	Dermal	Exposure		Primary	Ingestion	Inhalation	Dermal	Exposure
			<u> </u>				Routes Total	<u> </u>	Target Organ				Routes Total
Groundwater	Groundwater	Tap water	Acetone		<del>-</del> -	-	0.0E+00	Acetone	liver/kidney	1.9E-02	1.0E-01	2.7E-05	1.2E-01
			Trichloroethene	1.3E-07				Trichloroethene	_	1.9E-02	1.0E-01	7.5E-04	1.2E-01
			Tetrachloroethene	9.0E-07 1.8E-0 1.0E-07			1.2E-06	Tetrachioroethene	liver	1.7E-02	8.4E-03	2.0E-03	2.8E-02
		ļ	Chloroform	1.1E-08				Chloroform	liver	1.7E-03	1.1E+00	3.7E-05	1.1E+00
			Benzene	6.3E-08	1			Benzene	_	3.8E-03	3.6E-02	2.0E-04	4.0E-02
			Methylene chloride	7.2E-09	<b>8.2E-</b> 0	7.9E-11	1.5E-08	Methylene chloride	liv <b>er</b>	1.6E-04	6.0E-05	1.7E-06	2.2E-04
			(Total)	1.1E-06	1.5E-0	1.1E-07	2.7E-06	(Total)		6.1E-02	1.3E+00	3.0E-03	1.4E+00
				Total	Risk Acro: s(G	Proundweler)	2.7E-06	]	Total Hazard Index A	cross All Med	ia and All Expo	sure Routes	1.4E+00
			Total Risk	Across All Medi	ia and All i xpi	ceure Routes	2.7E-08	]			То	tal (liver) HI =	1.2E+00

NA = Not applicable.

#### TABLE 9.1.RME

SUMMARY OF FECEPTOR RISKS AND HAZARDS FOR COPCS

RE SONABLE MAXIMUM EXPOSURE

EVERGREEN MANC & SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timefram Current/Future Receptor Populatio Resident Receptor Age: Adult

Medium	Exposure Medium	Exposure Point	Chemical		C rosi	inagenic Risk		Chemical		Non-Carcin	ogenic Hazard	Quotient	:
				Ingestion	Inhala on	Dermal	Exposure		Primary	Ingestion	Inhalation	Dermal	Exposure
					L		Routes Total		Target Organ		_		Routes Total
Groundwater	Groundwater	Tap water	Acetone	-	-	-	•	Acetone	liver/kidney	2.7E-02	1.0E-01	1.0E-04	1.3E-01
1		1	Trichloroethene	6.2E-07	1.3E∹ ;	6.6E-08	2.0E-06	Trichloroethene	-	2.7E-02	1.0E-01	2.9E-03	1.3E-01
			Tetrachloroethene	4.4E-06	<b>6.3E</b> -∈7	1.4E-06	6.4E-06	Tetrachloroethene	liver	2.5E-02	8.4E-03	7.9E-03	4.1E-02
			Chloroform	5.2E-08	2.6E ∋	3.1E-09	2.6E-06	Chloroform	liver	2.5E-03	1.1E+00	1.5E-04	1.1E+00
			Benzene	3.1E-07	5.8E- :'	4.3E-08	9.3E-07	Benzene	-	5.5E-03	3.6E-02	7.7E-04	4.3E-02
			Methylene chloride	3.5E-08	2.8E- ∃	1.1E-09	6.4E-06	Methylene chloride	liver	2.3E-04	6.0E-05	6.9E-06	2.9E-04
			(Total)	5.4E-06	5.1E- 3	1.5E-06	1.2E-05	(Total)		8.8E-02	1.3E+00	1.2E-02	1.4E+00
				Total	Risk Acrc is(	Groundwater]	1.2E-05	,	Total Hazard Index A	cross All Med	ie and All Expo	sure Routes	1.4E+00
			Total Risk	Across All Medi	a and Ali Exp	posure Routes	1.2E-05	]			To	tal (liver) HI =	1.2E+00

NA = Not applicable.

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#### TABLE 9.2.CT

SUMMARY OF FECEPTOR RISKS AND HAZARDS FOR COPCS

C: NTRAL TENDENCY EXPOSURE

EVERGREEN MANO: SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timefram Current/Future Receptor Populatio Resident Receptor Age: Child

Medium	Exposure Medium	Exposure Point	Chemical		C⊾ ci	nogenic Riek		Chemical		Non-Carcin	ogenic Hazard	Quotient	
,			1	Ingestion	Inhelet: 1	Dermal	Exposure		Primary	Ingestion	Inhalation	Dermal	Exposure
							Routes Total		Target Organ				Routes Total
Groundwater	Groundwater	Tap water	Acetone				0.0E+00	Acetone	liver/kidney	4.5E-02	2.8E-01	4.5E-05	3.2E-01
		1	Trichloroethene	8.4E-08 2.9E-07 2.4E-09			3.7E-07	Trichloroethene	-	4.5E-02	2.8E-01	1.3E-03	3.2E-01
			Tetrachioroethene	6.0E-07 1.4E-C <b>5.1E-08</b>			7.9E-07	Tetrachioroethene	liver	4.0E-02	2.3E-02	3.4E-03	6.6E-02
		<b>,</b>	Chloroform	7.0E-09	5.8E-0	1.1E-10	5.9E-07	Chloroform	liver	4.0E-03	2.9E+00	6.4E-05	2.9E+00
			Benzene	4.2E-08	1.3E-0	1.6E-09	1.7E-07	Benzene	-	8.9E-03	9.8E-02	3.3E-04	1.1E-01
			Methylene chloride	4.6E-09	6.4E-0	3.8E-11	1.1E-08	Methylene chloride	liver	3.7E-04	1.6E-04	3.0E-06	5.4E-04
			(Total)	) 7.4E-07 1.1E-0 5.8E-08			1.9E-06	(Total) 1.4E-01 3.6E+00 5.2E-03			5.2E-03	3.7E+00	
	_		-	Total Risk Acro: :[Groundwater]			1.9E-06	Total Hazard Index Across All Media and All Exposure Routes				3.7E+00	
			Total Risk	Across All Medi	ia and All ι φ	ceure Routes	1.9E-06	Total (liver) HI ≃					3.3E+00

NA = Not applicable.

#### TABLE 9.2.RME

SUMMARY OF FECEPTOR RISKS AND HAZARDS FOR COPC®

RE. SONABLE MAXIMUM EXPOSURE

EVERGREEN MANC : SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timefram Current/Future Receptor Populatio Resident Receptor Age: Child

Medium	Exposure Medium	Exposure Point	Chemical		C. c	inogenic Risk		Chemical		Non-Carcin	ogenic Hazard	Quotient	
				Ingestion	Inhalat. n	Dermai	Exposure Roules Total		Primary Target Organ	Ingestion	Inhalation	Dermal	Exposure Roules Total
Groundwater	Groundwater	Tap water	Acetone	-		-	0.0E+00	Acetone	liver/kidney	6.4E-02	2.8E-01	1.8E-04	3.4E-01
			Trichloroethene	3.6E-07	3.6E-07 8.6E-4 2.9E-08			Trichloroethene	_	6.4E-02	2.8E-01	5.1E-03	3.5E-01
		İ	Tetrachioroethene	2.6E-06				Tetrachioroethene	liver	5.8E-02	2.3E-02	1.4E-02	9.4E-02
			Chloroform	3.0E-08				Chloroform	liver	5.8E-03	2.9E+00	2.5E-04	2.9E+00
	]	ļ	Benzene	1.8E-07	3.9E-€ ¹	1.9E-08	5.9E-07	Benzene	-	1.3E-02	9.8E-02	1.3E-03	1.1E-01
ļ			Methylene chloride	2.1E-08	1.9E-(	4.6E-10	4.0E-08	Methylene chloride	liver	5.3E-04	1.6E-04	1.2E-05	7.1E-04
			(Total)	3.2E-06	3.4E-( >	6.6E-07	7.3E-06	(Total)		2.0E-01	3.6E+00	2.1E-02	3.8E+00
	<del></del>			Total Risk Acro s[Groundwater]			7.3E-06	Total Hazard Index Across All Media and All Exposure Routes			3.8E+00		
			Total Risk	Across All Medi	ie and All 🗴	posure Routes	7.3E-06	Total (liver) HI = 3.4					3.4E+00

NA = Not applicable.

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#### TABLE 9.3.CT

#### SUMMARY OF FECEPTOR RISKS AND HAZARDS FOR COPCS

#### CI NTRAL TENDENCY EXPOSURE

EVERGREEN MANO ? SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Future Receptor Population: Commercial/Industrial Receptor Age: Adult

Medium	Exposure Medium	Exposure Point	Chemical		arcinog	jenic Risk		Chemical		Non-Carcin	ogenic Hazard	Quotient	
			•	Ingestion	In aletion	Dermal	Exposure		Primary	Ingestion	Inhalation	Dermal	Exposure
					L		Routes Total		Target Organ				Routes Total
Groundwater	Groundwater	Tap water	Acetone	-	F -	-	0.0E+00	Acetone	liver/kidney	6.8E-03	7.3E-02	1.9E-05	8.0E-02
			Trichloroethene	5.8E-08	: 4E-07	4.5E-09	4.0E-07	Trichloroethene	-	6.8E-03	7.3E-02	5.3E-04	8.1E-02
			Tetrachioroethene	4.1E-07 1 7E-07 9.6E-08			6.8E-07	Tetrachioroethene	liver	6.2E-03	6.0E-03	1.4E-03	1.4E-02
			Chloroform	4.8E-09				Chloroform	liver	6.2E-04	7.7E-01	2.7E-05	7.7E-01
			Benzene	2.9E-08	1 5 <b>E-07</b>	3.0E-09	1.9E-07	Benzene	-	1.4E-03	2.6E-02	1.4E-04	2.7E-02
		1	Methylene chloride	3.3E-09	7 5E-09	7.2E-11	1.1E-08	Methylene chloride	liver	5.7E-05	4.3E-05	1.2E-06	1.0E-04
					L			<b>]</b>					
			(Total)	5.1E-07	1 4E-06	1.0E-07	2.0E-06	(Total)		2.2E-02	9.5E-01	2.2E-03	9.7E-01
8.			-	Total Risk A rose[Groundwater]			2.0E-06	Total Hazard Index Across All Media and All Exposure Routes 9.76			9.7E-01		
			Total Ris	tisk Across All Media and III Exposure Routes			2.0E-06	Total (liver) HI = 8.6				8.6E-01	

N/A = Not applicable.

#### TABLE 9.3.RME

SUMMARY OF A ECEPTOR RISKS AND HAZARDS FOR COPCS

RE SONABLE MAXIMUM EXPOSURE

EVERGREEN MANC & SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenerio Timefreme: Future Receptor Population: Commerciel/Industrial Receptor Age: Adult

Medium	Exposure Medium	Exposure Point	Chemical		Carcinog	genic Risk		Chemical	Non-Carcinogenic Hazard Quotient					
			i	Ingestion	l: :alation	Dermal	Exposure		Primary	Ingestion	Inhalation	Dermal	Exposure	
							Routes Total		Target Organ				Routes Total	
Groundwater	Groundwater	Tap water	Acetone '	-	-	-	-	Acetone	liver/kidney	9.8E-03	7.3E-02	7.4E-05	8.3E-02	
		İ	Trichloroethene	2.3E-07	4E-07	4.9E-08	1.2E-06	Trichloroethene	-	9.8E-03	7.3E-02	2.1E-03	8.5E-02	
			Tetrachloroethene	1.6E-06 7E-07 1.0E-06			3.2E-06	Tetrachioroethene	liver	8.8E-03	6.0E-03	5.6E-03	2.0E-02	
			Chloroform	1.9E-06			1.9E-06	Chloroform	liver	8.8E-04	7.7E-01	1.0E-04	7.7E-01	
			Benzene	1.2E-07	3E-07	3.2E-08	5.8E-07	Benzene	-	2.0E-03	2.6E-02	5.5E-04	2.8E-02	
			Methylene chloride	1.3E-08	: 1E-06	7.9E-10	3.5E-08	Methylene chloride	liver	8.2E-05	4.3E-05	4.9E-06	1.3E-04	
			(Total)	2.0E-08	8E-06	1.1E-06	6.9E-06	(Total)		3.1E-02	9.5E-01	8.5E-03	9.9E-01	
			•	Total Risk / :rose[Groundwater]			6.9E-06	Total Hazard Index Across All Media and All Exposure Routes 9.9E-01					9.9E-01	
			Total Risk	Across All Media a	na Ji Expos	ure Roules	6.9E-06	Total (liver) Ht = 8.7				8.7E-01		

N/A = Not applicable.

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#### TABLE 10.1.CT

#### SUMMARY OF F ECEPTOR RISKS AND HAZARDS FOR COPCs

#### CI NTRAL TENDENCY EXPOSURE

EVERGREEN MANO SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timefram Current/Future Receptor Populatio Resident Receptor Age: Adult

Medium	Exposure Medium	Exposure Point	Chemical		C. cir	nogenic Riek		Chemical		Non-Carcin	ogenic Hazard	Quotient	
				Ingestion	inhalati n	Dermal	Exposure Routes Total		Primary Target Organ	Ingestion	Inhelation	Dermal	Exposure Routes Total
Groundwater	Groundwater	Tap water	Acetone	-	-	-	0.0E+00	Acetone	liver/kidney	1.9E-02	1.0E-01	2.7E-05	1.2E-01
			Trichloroethene	1.3E-07	3.7E-0	4.9E-09	5.0E-07	Trichloroethene	-	-	-	-	-
			Tetrachioroethene	9.0E-07	1 1 1			Tetrachloroethene	liver	1.7E-02	8.4E-03	2.0E-03	2 8E-02
,			Chloroform	1.1E-08	7.5E-0	2. <b>3E</b> -10	7.6E-07	Chloroform	liver	1.7E-03	1.1E+00	3.7E-05	1.1E+00
	ĺ	l	Benzene	6.3E-08	1.7E-0	3.2E-09	2.3E-07	Benzene	_	-	-	-	
			Methylene chloride	7.2E-09	8.2E-C	7.9E-11	1.5E-08	Methylene chloride	liver	1.6E-04	6.0E-05	1.7E-06	2.2E-04
			(Total)	1.1E-06	1.5E-€	1.1E-07	2.7E-06	(Total)		3.8E-02	1.2E+00	2.1E-03	1.2E+00
			-	Total Risk Acro (Groundwater)			2.7E-06	Total Hazard Index Across All Media and All Exposure Routes			1.2E+00		
			Total Risk	Across All Med	ia and All : xp	ceure Routes	2.7E-06	Total (liver) HI =					1.2E+00

NA = Not applicable.

#### TABLE 10.1.RME

#### SUMMARY OF I ECEPTOR RISKS AND HAZARDS FOR COPC#

RE .SONABLE MAXIMUM EXPOSURE

EVERGREEN MANC R SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timefram Current/Future Receptor Populatio Resident Receptor Age: Adult

Medium	Exposure Medium	Exposure Point	Chemical		C roin	nogenic Risk		Chemical		Non-Carcin	ogenic Hazard	Quotient	
				Ingestion	inheiat 10	Dermal	Exposure		Primary	Ingestion	Inhalation	Dermal	Exposure
Groundwater	Groundwater	Tap water	Acetone				Routes Total	Acetone	Target Organ liver/kidney	2.7E-02	1.0E-01	1.0E-04	Routes Total
			Trichloroethene	6.2E-07 1.3E-4 3 6.6E-08			2.0E-06	Trichloroethene	-	-	-	-	-
ļ			Tetrachloroethene	4.4E-06 6.3E-4 / 1.4E-06			6.4E-06	Tetrachioroethene	liver	2.5E-02	8.4E-03	7.9E-03	4.1E-02
1	!		Chloroform	5.2E-08	2.6E⊣ ;	3.1E-09	2.6E-06	Chloroform	liver	2.5E-03	1.1E+00	1.5E-04	1.1E+00
			Benzene	3.1E-07	5.8E⊣ ′	4.3E-08	9.3E-07	Benzene	-	-	-	- 1	-
			Methylene chloride	3.5E-08	2.8E-⊢ ∃	1.1E-09	6.4E-08	Methylene chloride	liver	2.3E-04	6.0E-05	6.9E-06	2.9E-04
			(Total)	5.4E-06	5.1E⊰ ∋	1.5E-06	1.2E-05	(Total)		5.5E-02	1.2E+00	6.2E-03	1.2E+00
•	_		<del>-</del>	Total Risk Acrc s[Groundwater]			1.2E-05	Total Hazard Index Across All Media and All Exposure Routes 1.2				1.2E+00	
			Total Risk	Across All Media and All I xxxxxxx Routes			1.2F-05	Total (liver) HI ≈ 1.25-					1.2E+00

NA = Not applicable.

#### TABLE 10.2.CT

#### SUMMARY OF ECEPTOR RISKS AND HAZARDS FOR COPCS

C. NTRAL TENDENCY EXPOSURE

EVERGREEN MANC & SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timefram Current/Future Receptor Populatio Resident Receptor Age: Child

Medium	Exposure Medium	Exposure Point	Chemical		Circin	nogenic Riek		Chemical	Non-Carcinogenic Hazard Quotient					
				Ingestion	inhelat in	Dermat	Exposure Routes Total		Primary Target Organ	Ingestion	Inhalation	Dermal	Exposure Routes Total	
Groundwater	Groundwater	Tap water	Acetone	-	-	-	0.0E+00	Acetone	liver/kidney	4.5E-02	2.8E-01	4.5E-05	3.2E-01	
			Trichtoroethene	8.4E-08	2.9E-€ /	2.4E-09	3.7E-07	Trichloroethene	-	-	-	-	-	
			Tetrachloroethene	6.0E-07 1.4E-4 5.1E-08			7.9E-07	Tetrachioroethene	liver	4.0E-02	2.3E-02	3.4E-03	6.6E-02	
ŀ	Ī		Chloroform	7.0E-09	5.8E≺ '	1.1E-10	5.9E-07	Chloroform	liver	4.0E-03	2.9E+00	6.4E-05	2.9E+00	
			Benzene	4.2E-08	1.3E-(	1.6E-09	1.7E-07	Benzene	_	-	-	-	-	
1			Methylene chloride	4.8E-09	6.4€-( )	3.8E-11	1.1E-08	Methylene chloride	liver	3.7E-04	1.6E-04	3.0E-06	5.4E-04	
			(Total)	7.4E-07	1.1E-( i	5.5E-08	1.9E-06	. (Total)		8.9E-02	3.2E+00	3.5E-03	3.3E+00	
	· · · · · ·	3.5	<del></del>	Total	Risk Acrc s[G	Proundweler]	1.9E-06	Total Hazard Index Across All Media and All Exposure Routes 3.				3.3E+00		
			Total Rick	Across All Med	ie end Att -vm	onera Routes	1.9F-06	Total (liver) HI = 3.3E+0					3.3E+00	

NA = Not applicable.

#### TABLE 10.2.RME

SUMMARY OF I ECEPTOR RISKS AND HAZARDS FOR COPC#

RE SONABLE MAXIMUM EXPOSURE

EVERGREEN MANC SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timefram Current/Future Receptor Populetio Resident Receptor Age: Child

Medium	Exposure Medium	Exposure Point	Chemical		<b>C</b> . •	cinogenic Risk		Chemical		Non-Carcin	ogenic Hazerd	Quotient	
1			i	Ingestion	inhalat r	n <b>Dermal</b>	Exposure	1	Primary	Ingestion	Inhelation	Dermal	Exposure
							Routes Total		Target Organ				Routes Total
Groundwater	Groundwater	Tap water	Acetone	_	-	T -	0.0E+00	Acetone	liver/kidney	6.4E-02	2.8E-01	1.8E-04	3.4E-01
ł	l	ł	Trichloroethene	3.6E-07	8.6E-€	2.9E-08	1.2E-06	Trichloroethene	_	-	-	-	-
1			Tetrachioroethene	2.6E-06	4.3E-{	6.1E-07	3.6E-08	Tetrachioroethene	liver	5.8E-02	2.3E-02	1.4E-02	9.4E-02
		ļ	Chloroform	3.0E-08	1.7E-C	1.3E-09	1.8E-06	Chloroform	liver	5.8E-03	2.9E+00	2.5E-04	2.9E+00
		ŀ	Benzene	1.8E-07	3.9E-(	1.9E-08	5.9E-07	Benzene	-	-			-
			Methylene chloride	2.1E-08	1.9E-C	4.6E-10	4.0E-08	Methylene chloride	liver	5.3E-04	1.6E-04	1.2E-05	7.1E-04
			(Total)	3.2E-06	3.4E-(	6.6E-07	7.3E-06	(Total)		1.3E-01	3.2E+00	1.4E-02	3.4E+00
				Total	Riek Acro	[Groundwater]	7.3E-06	Total Hazard Index Across All Media and All Exposure Routes 3.4					3.4E+00
			Total Risk	Across All Medi	ia and All	posure Routes	7.3E-06	Total (liver) HI = 3					3.4E+00

NA = Not applicable.

#### TABLE 10.3.CT

SUMMARY OF # ECEPTOR RISKS AND HAZARDS FOR COPCs

C: NTRAL TENDENCY EXPOSURE

EVERGREEN MANO:: SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Future Receptor Population: Commercial/industrial Receptor Age: Adult

Medium	Exposure Medium	Exposure Point	Chemical		arcinog	enic Riek		Chemical		Non-Carcinogenic Hazard Quotient				
				Ingestion	Ir aletion	Dermai	Exposure		Primary	Ingestion	Inhalation	Dermal	Exposure	
							Routes Total	<u></u>	Target Organ				Routes Total	
Groundwater	Groundwater	Tap water	Acetone	-		-	0.0E+00	Acetone	liver/kidney	-	-	-	-	
			Trichloroethene	5.8E-08 4E-07 4.5E-09			4.0E-07	Trichloroethene		-	-	-	-	
			Tetrachloroethene	4.1E-07 7E-07 9.8E-08			6.6E-07	Tetrachloroethene	liver	-	-	-	-	
			Chloroform	4.1E-07   7.E-07   9.8E-08   4.8E-09   (.9E-07   2.1E-10			6.9E-07	Chloroform	liver	-	-	-	- 1	
			Benzene	2.9E-08	- 5 <b>E-0</b> 7	3.0E-09	1.9E-07	Benzene			-	-	- 1	
			Methylene chloride	3.3E-09	7 5 <b>E-09</b>	7.2E-11	1.1E-08	Methylene chloride	liver	-			-	
							•					·······		
			(Total)	5.1E-07	4E-06	1.0E-07	2.0E-06	(Total)			_			
				Total Ris	k A ross[Gr	zundwaler)	2.0E-06	Total Hazard Index Across All Media and All Exposure Routes					_	
			Total Risi	sk Across All Media and Il Exposure Roules			2.0E-06	Total (liver) HI =						

N/A = Not applicable.

RFW036-2A-AHV RFW036-2A-AGQP

#### TABLE 10.3.RME

SUMMARY OF A ECEPTOR RISKS AND HAZARDS FOR COPCS

RE SONABLE MAXIMUM EXPOSURE

EVERGREEN MANO R SITE, ROSCOE, WINNEBAGO COUNTY, ILLINOIS

Scenario Timeframe: Future Receptor Population: Commercial/Industrial Receptor Age: Adult

Medium	Exposure Medium	Exposure Point	Chemical		Carcino	genic Risk		Chemical		Non-Carcin	ogenic Hezard	Quotient	
				Ingestion	Ir aletion	Dermal	Exposure	Ì	Primary	Ingestion	Inhelation	Dermal	Exposure
						L	Routes Total		Target Organ				Routes Total
Groundwater	Groundwater	Tap water	Acetone	-	-	-	-	Acetone	liver/kidney	_	-		-
			Trichloroethene	2.3E-07	: 4E-07	4.9E-08	1.2E-06	Trichloroethene	_	-	-	-	-
			Tetrachloroethene	1.6E-06	7E-07	1.0E-06	3.2E-06	Tetrachioroethene	liver		-	-	-
			Chloroform	1.9E-08	9E-06	2.3E-09	1.9E-06	Chloroform	liver	-	-	-	-
			Benzene	1.2E-07	3E-07	3.2E-08	5.8E-07	Benzene	-	-	-	] -	-
			Methylene chloride	1.3E-08	: 1E-08	7.9E-10	3.5E-08	Methylene chloride	liver	-	-	-	-
			(Total)	2.0E-06	8E-06	1.1E-06	6.9E-08	(Total)			-	-	-
				Total Risk / ross[Groundwater]			6.9E-06	Total Hazard Index Across All Media and All Exposure Routes -				-	
			Total Ris	k Across All Media a	nd Jil Expos	ure Roules	6.9E-08	Total (liver) HI ≖				-	

N/A = Not applicable.

RFW036-2A-AHV RFW036-2A-AGQP

## APPENDIX B

FRACTURE TRACE ANALYSIS, RESOLUTION RESOURCES, INC.

## LETTER REPORT

## FRACTURE TRACE ANALYSIS

### at the

# EVERGREEN MANOR SITE Winnebago County, Illinois

Submitted to:

ROY F. WESTON, INC.

3 Hawthorne Parkway, Suite 400
Vernon Hills, Illinois 60061
Mr. Kurt T. Fischer
Tel: 847-918-4016 • Fax: 847-918-4055

Submitted by:

Resolution Resources, Inc.

Brian B. Herridge 310 West 52nd Street Minneapolis, MN 55419 612-824-3234

Mary-Linda Adams 8167 Old Waterloo Road Warrenton, VA 22186 540-349-9176

## TABLE OF CONTENTS

I.	PR	ROJECT OVERVIEW	1
	1.1	INTRODUCTION	1
	1.2	OBJECTIVES	1
,	1.3		
	1.3.		
	1.3.		
	1.3.		
2.	BA	CKGROUND INFORMATION	2
2	2.1	LOCATION AND HISTORY	2
	2.2	SITE GEOLOGY & HYDROLOGY	
•	nr.		
<b>J</b> .	PH	IOTOGRAPHIC ANALYSIS	د
3	3.1	LINEAMENTS	5
3	3.2	HISTORICAL ANALYSIS	6
3	3.3	FRACTURE TRACE ANALYSIS	7
4.	RE	COMMENDATIONS	8
4	.1	SAMPLE POINT LOCATIONS	8
4	.2	HISTORICAL PHOTOGRAPHS	
5.	DIS	SCUSSION	9
6.	RE	FERENCES	10

## LIST OF TABLES, FIGURES & APPENDICES

Figure 1	Topographic Map			
Figure 2	Site Plan, Sampling Locations, & VOC Plume			
Figure 3	Fault Blocks and Lineaments Schematic			
Figure 4	Fractured Column Schematic			
Figure 5	1939 Historical Photographs			
Figure 6	1946 Historical Photograph			
Figure 7	1970 Historical Photographs			
Figure 8	1978 Historical Photographs			
Figure 9	1994 Historical Photograph			
Figure 10	Fracture Trace Analysis: 1978 Aerial Photographs			
Figure 11	Fracture Trace Analysis & Site Plan Composite			
Figure 12	Recommended Sample Point Locations			
Γable 1	Available Aerial Photographs of the Site			
Гable 2	Examined Aerial Photographs			
Appendix A	Uninterpreted Aerial Photographs			
Figure A1	1978 Aerial Photograph: 1-4-1			
Figure A2	1978 Aeriai enotograpii. 1-4-2			
Figure A3	1978 Aerial Photograph: 2-1-1			
Figure A4	1978 Aerial Photograph: 2-1-2			
Figure A5	1978 Aerial Photograph: 2-1-3			
Figure A6	1978 Aerial Photograph: 2-1-4			
Figure A7	1978 Aerial Photograph: 2-1-5			

## 1. PROJECT OVERVIEW

#### 1.1 INTRODUCTION

Resolution Resources, Inc. (RRI) performed a fracture trace analysis on aerial photographs of the Evergreen Manor Site, located in Winnebago County, Illinois. Figure 1 shows the site on the South Beloit, Illinois-Wisconsin topographic map (USGS, 1993). The site area is located southeast of Rockton in the relatively flat lying valley of the Rock River. The site includes four residential subdivisions that were developed from 1940 to 1988 on farmland. It is bounded by the Rock River to the south and surrounded by forest and farmland. EPA sampling results from 1993 and 1994 indicated that TCA and TCE had impacted over 60 residential wells. It has been estimated that contaminants have affected 700 people in 250 residences.

### 1.2 OBJECTIVES

The objectives of the photographic interpretation at the Evergreen Manor site were to help select representative groundwater sample locations, and to further evaluate the suspected source areas and the extent of the VOC plume. This information will be used to focus the investigation and to help evaluate the remedial alternatives at the site. The final report includes a series of suggested sample points to better define the plume. These points have been plotted on the aerial photographs. A discussion for the location of each point has been included in the report.

#### 1.3 WORK TASKS

In order to meet the objectives of the work, the following tasks were performed:

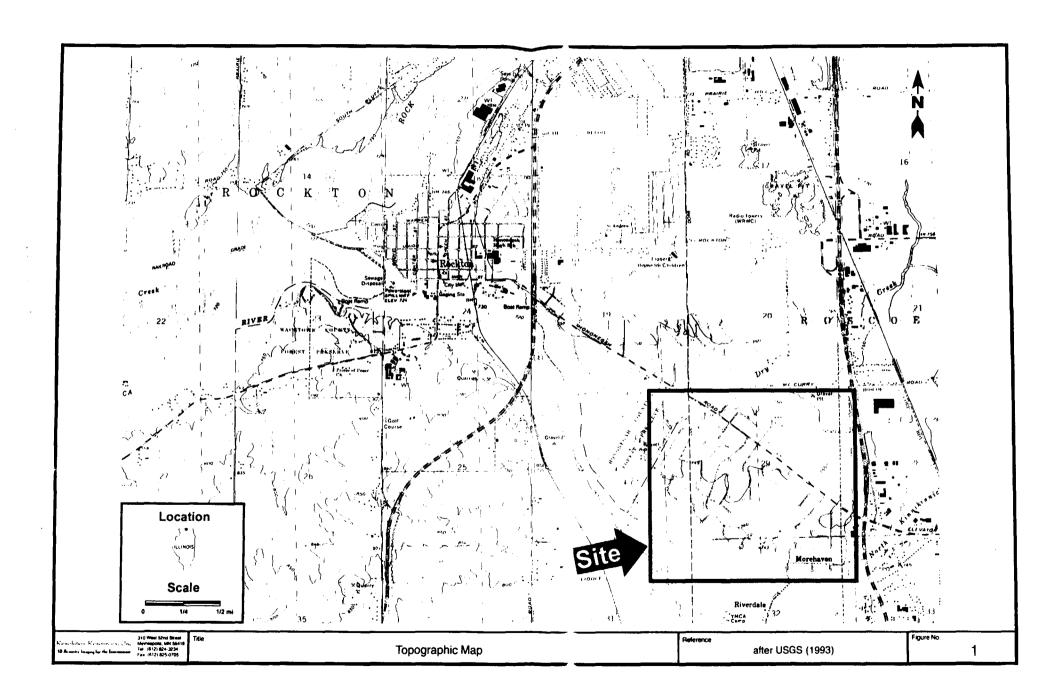
Task 1 Background Review

Task 2 Photographic Interpretation

Task 3 Report

#### 1.3.1 BACKGROUND REVIEW

This task included reviewing available background information on the site history, contaminant distribution, geology, and hydrogeology, which was provided by Roy F. Weston, Inc. The task also included performing a search for and the selection of the most appropriate historical aerial photographs.



## 1.3.2 PHOTOGRAPHIC INTERPRETATION

This task consisted of reviewing available aerial photographs and then performing a stereographic interpretation. Selected photographs were evaluated using a Topcon viewer. Historical photographs were also reviewed to better discuss possible source locations.

#### 1.3.3 REPORT

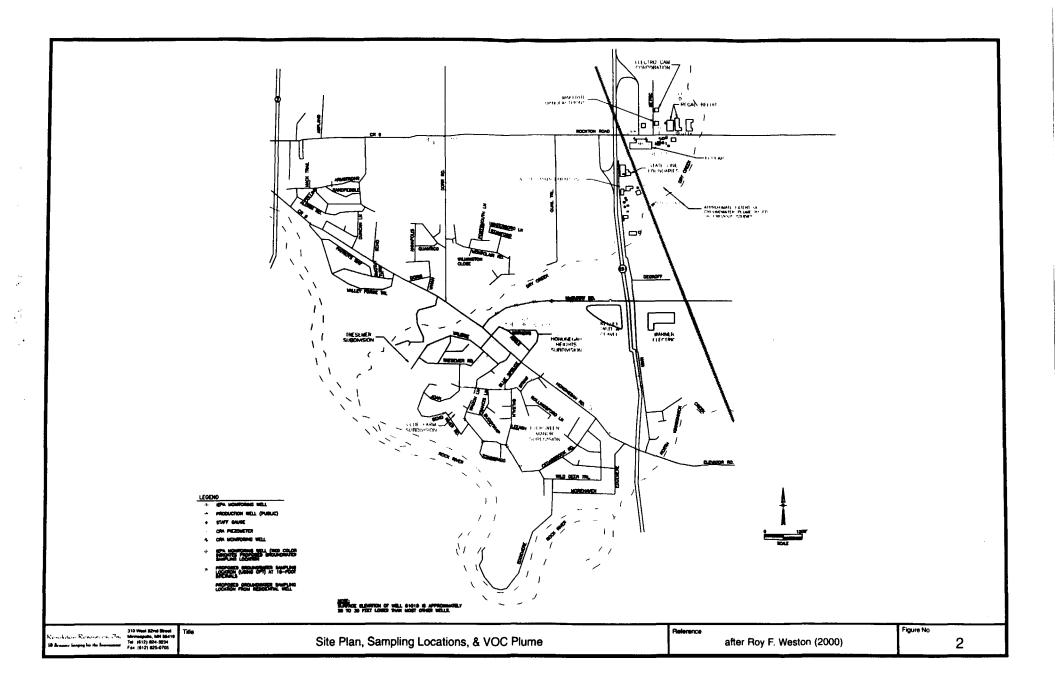
This letter report summarizes all activities that were performed on the project. A hard copy of the photographic interpretation has been included, along with recommendations for the location of sample points.

## 2. BACKGROUND INFORMATION

### 2.1 LOCATION AND HISTORY

The Evergreen Manor site is located in Winnebago County, Illinois about 1.5 miles northwest of Roscoe. The site has been defined by the extent of a solvent plume in groundwater. It includes four residential subdivisions, as shown on Figure 2: Hononegah Heights, developed between 1940 and 1964; Tresemer, developed between 1972 and 1974; Olde Farm, developed between 1976 and 1979; and Evergreen Manor, developed between 1986 and 1988. Hononegah Country Estates subdivision, which is characterized by a separate contaminant plume, is located further to the southeast. Prior to development the land was farmed. As shown on the topographic map (Figure 1), the site is bounded to the south by the Rock River, by the Hononegah Forest Preserve to the west, and farmland to the east and north. Kelly Sand and Gravel, a gravel pit and concrete mixing facility, is located about ½ mile northeast of the site. Further to the northeast, about two miles, is an industrial park.

The presence of VOCs in the groundwater at Evergreen Manor was first detected in the 1990's as a result of requirements of a lending institution. Further sampling showed that a VOC plume consisting of TCE; 1,1 DCE; C-1,2 DCE; 1,1 DCA; TCA; 1,12 TCA; and PCE existed beneath Hononegah Heights and Evergreen Manor. The site was added to CERCLIS in 1991, and a Preliminary Assessment (PA) was performed in 1992. A Sampling Site Inspection (SSI) was performed in 1992 to gather further information for the Hazard Ranking System (HRS). Soil gas and groundwater samples were collected. No contaminants were found north of Rockton Road. VOCs were detected in a well north of the Waste Management facility (Figure 2).



Based on the previous results, an Expanded Site Inspection (ESI) was conducted in 1993, which consisted of the collection of 49 water samples from 45 residential wells in the Hononegah Heights, Olde Farm and Evergreen Manor Subdivisions. By 1994 a total of 267 drinking water wells were sampled, with 108 wells above the MCL and 203 that showed VOCs. A total of 24 monitor wells have also been installed and tested. The source area has been determined to be located near the intersection of Rockton Road and Route 251. Four PRPs have been identified and have declined to participate in the remedial effort (Weston, 1999).

## 2.2 SITE GEOLOGY & HYDROLOGY

The site is located within the preglacial Rock River Valley, which is a bedrock valley that has been infilled with Quaternary glacial deposits. Highly permeable sand and gravel are the principal sediments. These deposits can reach thickness' of up to 300 feet. Logs from the wells drilled at the site, as deep as 100 feet, are characterized by sand and gravel deposits. The Rock River has eroded into Ordovician and Cambrian clastic and carbonate rocks, which were deposited on Precambrian granite.

The aquifers beneath the site consist of the glacial outwash sand and gravel and the St. Peter, Ironton-Galesville, and Mt. Simon Formation sandstones. Most of the domestic wells are drilled in the glacial deposits, from 50 to 80 feet, while larger municipal wells extend into the sandstone. The water table is about 35 feet below ground surface (bgs). The dolomite (carbonate) acts as an aquitard. Groundwater flow in the dolomite is through vertical fractures. Although the yields are not as high as in the glacial deposits or the sandstone some water wells have been drilled within the dolomite.

## 3. PHOTOGRAPHIC ANALYSIS

Aerial photography and satellite imagery can often provide very useful information on site history, including contaminant source area location, site development over time, regional topography, possible impacts to the local watershed, changes in surrounding commercial development, possible offsite contaminant sources and other influences and effects such as stressed vegetation. More importantly, stereographic photographs are an invaluable tool for the identification of geologic structural features such as fractures, faults and relative highs and lows (structural/topographic).

A library search was made to identify the years and scales of photographs that were available for the site. Table 1 lists all the photographs that are available. Table 2 references the date and scale for each photograph that was ordered and reviewed. Stereographic pairs of aerial photographs of the Evergreen Manor Site were examined for suitability to perform a fracture trace analysis.

TABLE 1: AVAILABLE AERIAL PHOTOGRAPHS OF THE SITE

Flight Year	Scale
1939	1" = 1667'
1945	1" = 2257'
1946	1" = 2267'
1951-53	1" = 1667'
1958	1" = 1667', 1" = 5500'
1964	1" = 1667'
1966	1" = 1000'
1970	1" = 1667', 1" = 1700', 1" = 3166'
1975	1" = 3167'
1976	1" = 3000'
1978	1" = 900', 1" = 2000'
1979	1" = 3300'
1980-81	1" = 4800', 1" = 6700'
1981	1" = 1000'
1986-88	1" = 1000', 1" = 3300', 1" = 3333', 1" = 4800', 1" = 5700'
1988	1" = 3300'
1991	1" = 2083'
1992-93	1" = 3300'
1994	1 <sub>n</sub> = 1000,

**TABLE 2: EXAMINED AERIAL PHOTOGRAPHS** 

Flight Year	Scale	Number
1939	1" = 1667'	4-25, 4-26, 4-38, 4-39
1946	1" = 2267'	852 - 854, 861 - 863
1970	1" = 1700'	2-150, 2-151, 2-152, 2-160, 2-161, 2-162
1978	1" = 900'	2-1-1 to 2-1-5, 1-4-1 to 1-4-2
1994	1" = 3600'	5775-92, 5775-93, 5775-94

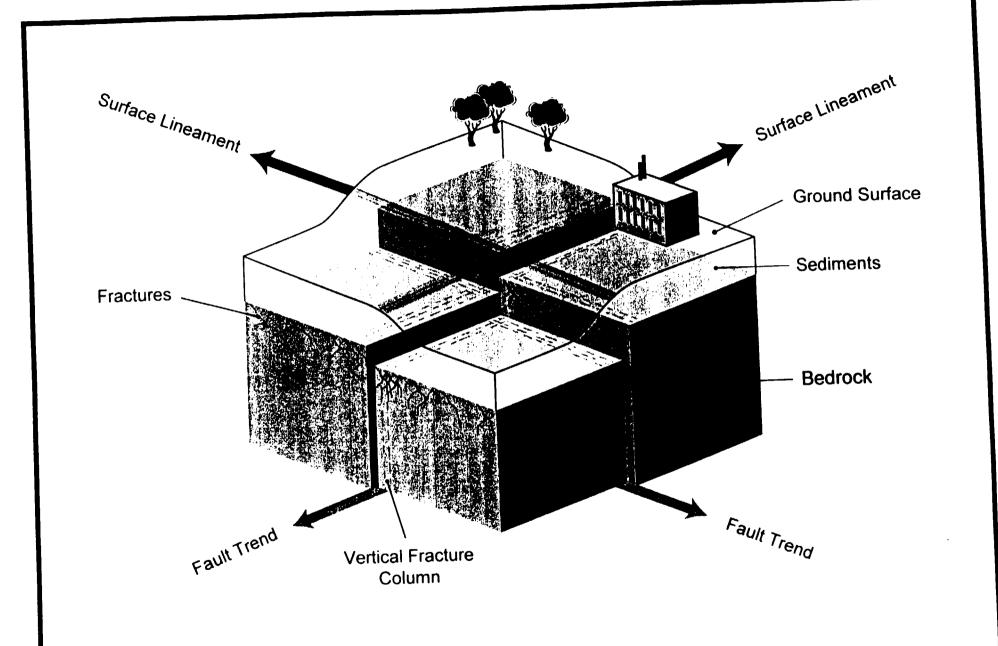
#### 3.1 LINEAMENTS

It is often the case that fractures and faults in bedrock or basement rock are propagated up through unconsolidated sediments to the surface as failure planes. This may be a result of occasional seismic activity and water movement along the fracture. The surface features caused by the fractures are called lineaments. Lineaments show subtle surface expressions that reveal their subsurface existence and can often be seen in aerial photographs. Lineaments are identified primarily based upon subtle changes in the shading and in the topography at the ground surface. Lineaments cut across different surface terrain and often display a topographic expression where one side of the lineament is slightly higher than the other side, as though offset has occurred. Materials that infill faults or fractures frequently have a different shading than the surrounding surfaces which have never been fractured.

Investigators (Culbreth, 1988; Wobber, 1967; Parizek, 1977; Rumsey, 1971) have found that lineaments can be identified in aerial photographs, even when sediments overlie the bedrock for hundreds of feet, and that they are manifestations of fractures or faults that have been propagated from bedrock to the surface through unconsolidated sediments and soil. The lineaments can be expressed by a variety of features (Hough, 1960), such as tonal changes in soil, changes in the directions of streams, straight segments in drainage patterns, or alignment of vegetation (since fractures are often more permeable, more water is available for enhanced growth of the plants). As a result of their work on LandSat imagery, compared to outcrop patterns and geophysical data in Montana and Wyoming, Marrs and Rains (1984) concluded that the lineaments represented the surface expression of boundaries of crustal blocks that have been activated throughout time.

Figure 3 shows a schematic diagram of how vertical fractures in the bedrock can propagate upward, and are expressed as surface lineaments. Since the major fracture systems have often been active through time (occasional reactivation), the sediments above bedrock have failure planes that are localized zones of weakness along which channels may cut. This explains why channels below the ground surface, but above the bedrock, can often be identified through the careful analysis of lineaments.

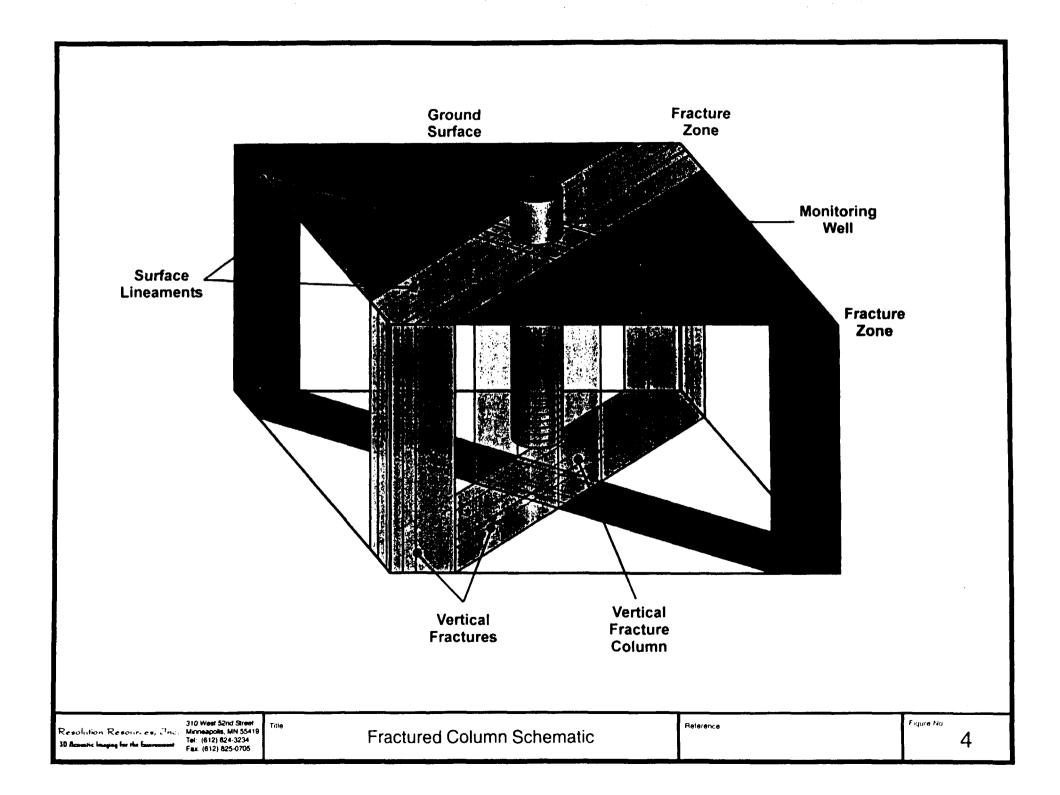
Fractures can have an effect on a variety of site characteristics including groundwater flow, contaminant transport, and well yield. Figure 4 schematically shows how the juncture of two vertical fractures can provide the optimum position for a vertical well screen, and if in connection with the source, for contaminant recovery. The vertical fractured column, created at the intersection of two vertical fracture sets, is connected to the other sub-vertical and horizontal fractures in the subsurface, making it the ideal location for a vertical well screen. Identifying fracture sets is very important to developing accurate site assessments and in designing effective remedial systems.



Resolution Resources, The Minnapolis, MN 55419
Tel: (612) 824-3234
Fax: (612) 825-0705

Fault B. . . ks and Lineaments Schematic (

Reference



### 3.2 HISTORICAL ANALYSIS

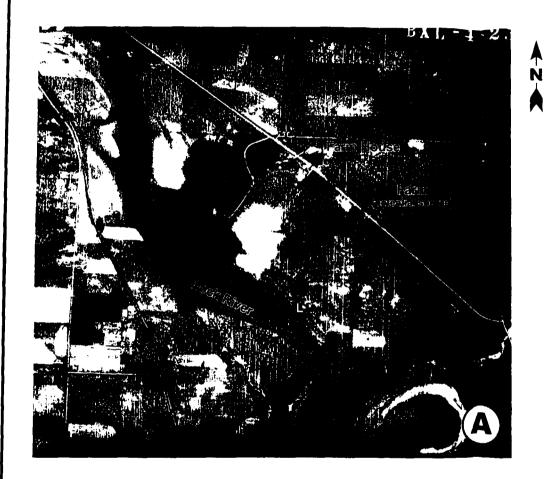
Photographs from 1939 to 1994 were analyzed to examine how the site has changed over the 55 year span that photographs are available. Figure 5a shows the area that is now occupied by the subdivisions in the 1939 vintage photographs. These photographs were examined because they predate development of the Hononegah Heights Subdivision, when the site was still devoted to farming activities. The quality of the photographs are the least of all the years that were examined, and a light dusting of snow, which was present when the photographs were taken, further obscures some features. Despite the grainy nature of the photographs, it is possible to see a low area, marked by an "L", to the east of a lineament. Lineaments are shown in red on the historical photographs, and dotted where inferred. It is reasonable to assume that ground and surface water would have drained to this low. A farmhouse and surrounding buildings are located south of Hononegah Road. An area circled in blue shows what is most likely a large farm. This area may be of interest for sampling, since solvents could have been used on farm vehicles. Figure 5b shows the suspected source area (shown by a blue circle), 2 miles to the northeast of the site, as it appeared in 1939. The area is completely rural with a farm and a house.

Figure 6 shows the site, including the area where the subdivisions were eventually located, and the suspected source areas in 1946. According to the background information construction on Hononegah Heights began in 1940, however there is little evidence from the 1946 photographs that any homes had been constructed. Two farms with outlying buildings, which are seen in the 1939 photographs, are still present in the subdivision area. The low at the river, seen on the 1939 photographs, is still present. Two new homes or farms have been added, since 1939, on either side of Hononegah Road. The quarry operations have begun by this date, and a farmhouse or office building is located near the quarry.

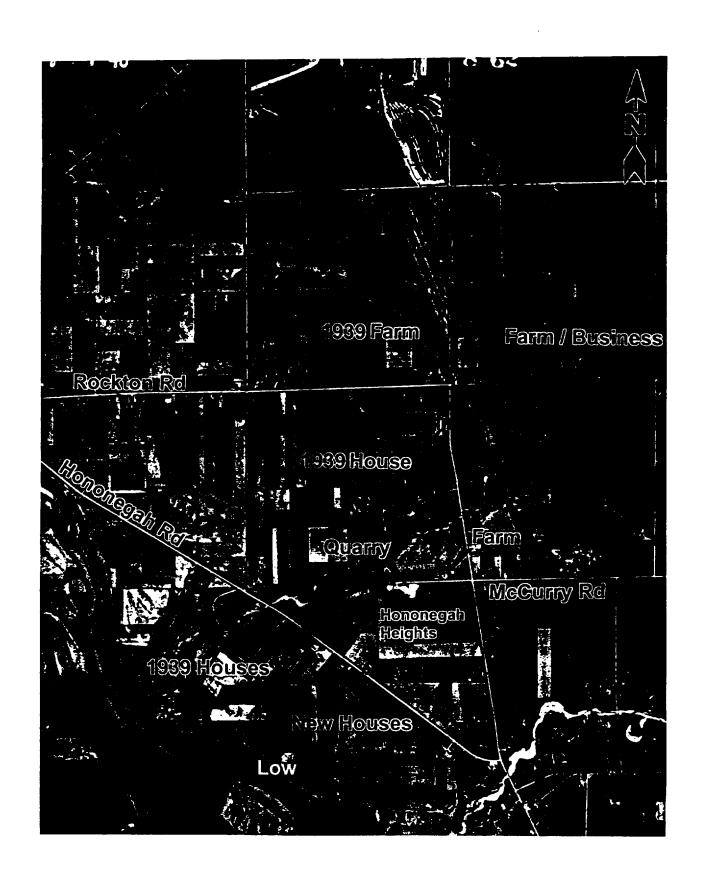
In the suspected source area, two homes or farms present in 1939 are still in existence, but a new farm or business has been added north of Rockton Road.

Figure 7a shows the subdivision area in 1970. The low by the river is still apparent, and a northeast trending lineament is prominent. The two houses present from 1939 are still in existence, as are two houses seen for the first time in the 1946 photographs. The Hononegah Heights Subdivision has clearly been developed, as has the quarry. An apparent upthrown block (labeled with a "U"), along an east-west trending fracture, is also present. The downthrown side is labeled with a "D". This upthrown block could act as a barrier to groundwater flow, and wells or piezometers should be constructed on either side of this block to determine if water levels change across the block. However, as a result of the high permeability of the sediments in the area a difference in hydraulic head across the upthrown block may not be observed.

Figure 7b shows the suspected source area in the 1970 set of photographs. The two farms/houses that were located west of Route 251 in the 1939 photographs are no longer







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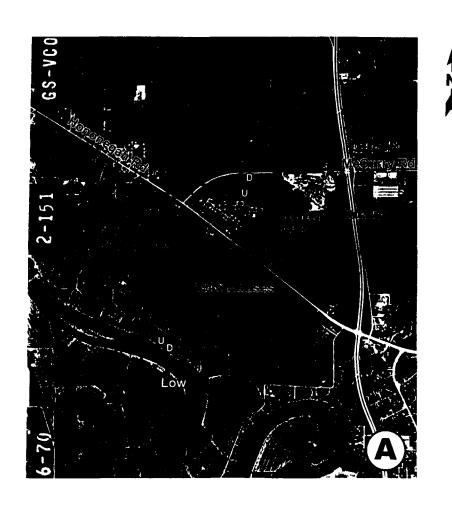
310 West 52nd Street Minneapolis, MN 55419 Tel. (612) 824-3234 Fax: (612) 825-0705

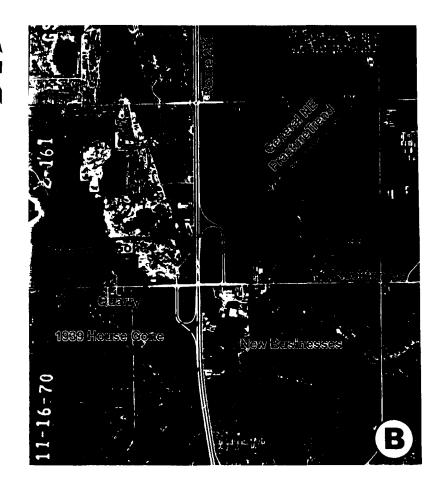
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1946 Historical Photograph

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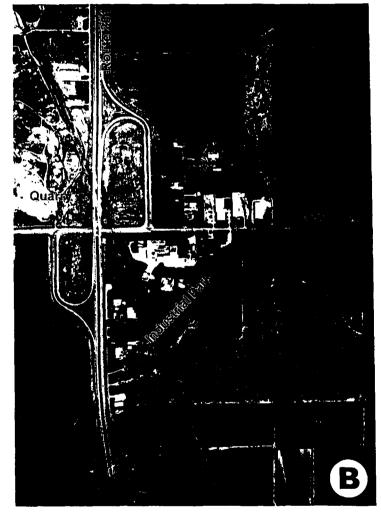
1970 Historical Photographs

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7





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1978 Historical Photographs

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1994 Historical Photograph

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there. Quarry operations have commenced. A farm to the east of Route 251 seen in 1946 is gone, and three new businesses have been constructed. A general northeast trend in lineaments, which is also observed in the dry streambed, is shown in the photographs.

Figure 8 shows the 1978 photographs. Figure 8a shows the extent of development in the subdivision area, and a potential source location from the 1939 photographs. Figure 8b shows the growth of businesses in the suspected source area east of Route 251, both south and north of Rockton Road.

Figure 9 shows the site in 1994, which is the most recent photograph that was obtained This photograph was taken at a regional scale so it is possible to see the general north-south and east-west, and northeast-southwest and northwest-southeast, linear trends present in the area

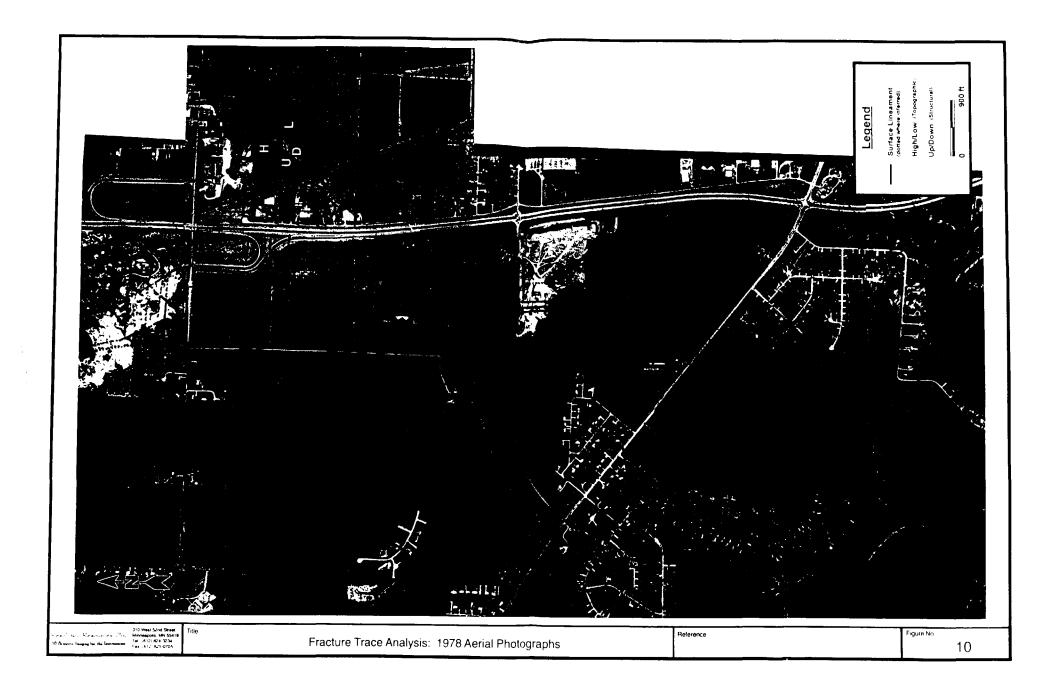
### 3.3 FRACTURE TRACE ANALYSIS

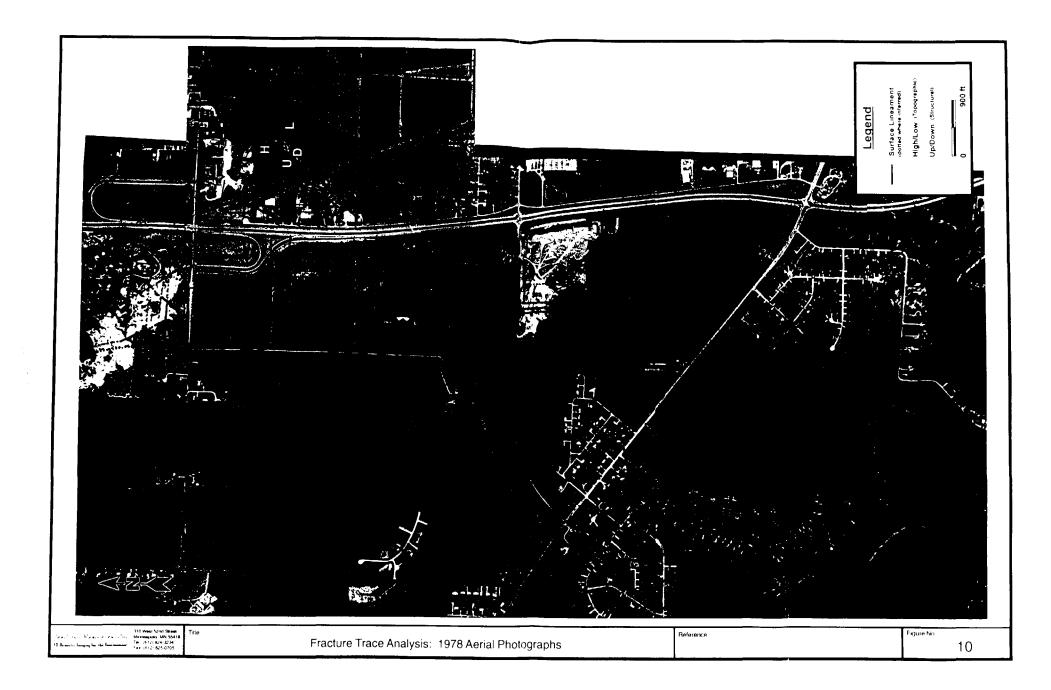
All of the photographs were evaluated for sun angle, seasonal variations, and greatest vertical exaggeration. The photographs from 1978 were chosen to perform a detailed fracture trace analysis, because of their scale and clarity. The uninterpreted, 1978 stereographic aerial photographs are included in **Appendix A**.

Figure 10 shows the mosaic of the detailed fracture trace analysis, which was performed on the 1978 photographs. Lineaments are shown in red, and are dotted where inferred. The fracture trace analysis shows that there are two main sets of fracture trends, a north-south/east-west set, and a northeast-northwest conjugate set. These two sets of lineaments are also seen in trends in the geomorphology. The clearest linear features can be seen in the quarry north of Rockton Road. There are many lineaments that trend northeast toward the subdivision from the industrial park. So a more permeable transport of fluids along fractures does exist. Once south of Hononegah Road the plume appears to follow north-south oriented fractures that extend north to the Rock River.

Figure 11 shows the lineaments from the fracture trace analysis superimposed with the site plan. Since the examined photographs were not orthorectified to remove distortions, the fit of the photographic interpretation on the site plan cannot be exact. The position of the lineaments on the site diagram should be regarded as a close approximation.

A cursory examination of the figure shows that in the area of the subdivisions, where the plume moves in a more north-south direction south of Hononegah Road, the most prominent lineaments are also oriented north-south. In the area north and northeast of Hononegah Road the most prominent lineaments trend northeast in the same direction as the plume. As an aside, northwest trending lineaments are most prominent to the north of the dry creek bed. There is some evidence then, that the direction of the VOC plume is controlled by fractures.





## 4. RECOMMENDATIONS

### 4.1 SAMPLE POINT LOCATIONS

It is necessary to mention that one of the limitations of aerial interpretation of fractures is plowed fields, which are in abundance at this site. It is impossible in many cases to see lineaments in freshly plowed fields. As a result, some of the sample locations are restricted in these areas.

Figure 12 shows the recommended locations for a total of 44 sample points (SP) on fractures. Locations 1-9 are at the juncture of one or more fractures or in lows in the southernmost part of the plume. Sample points 3 and 5 are in lows. It will be very important to try to place the actual sample points as close to the recommended locations as is possible, since these more permeable fracture zones are quite narrow.

SP 10 is at the juncture of a north and northeast trending fracture. Note that northeast trending Lineaments A and B to the east may coincide with the high yields necessary for public production wells, which are noted on Figure 2.

SP 11 and 13 are located along northeast trending fractures that extend toward the suspected source areas

SP 12 and 13 are at the juncture of a north-south trending fracture and the northeast trending fractures.

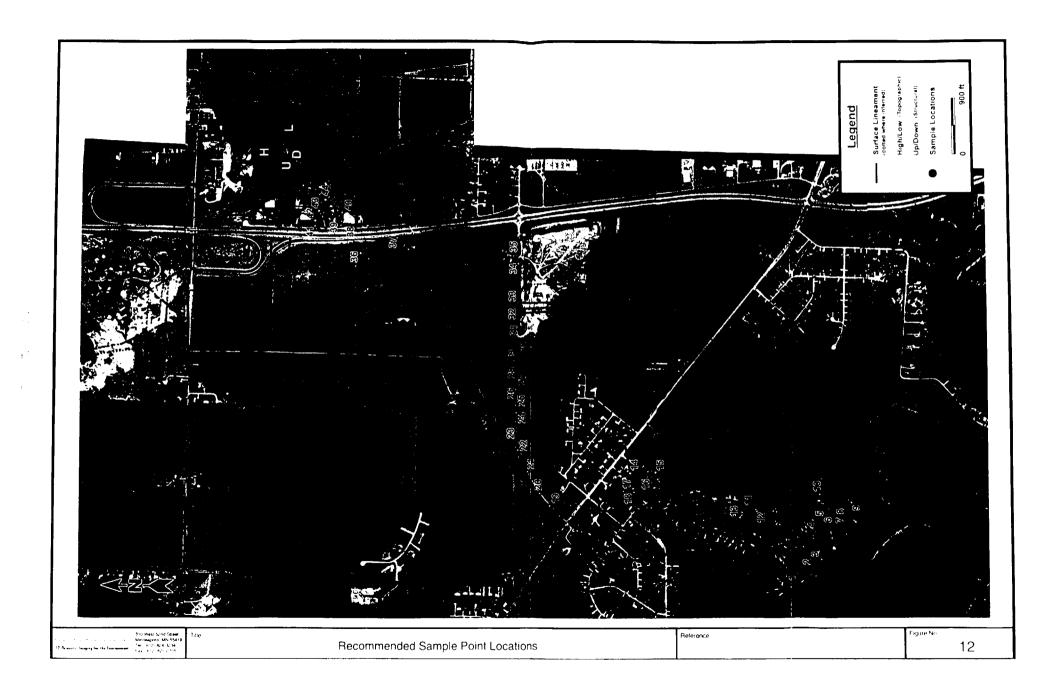
SP 14-18 are located along fractures down-gradient of the large farm, which has been identified from the 1939 photographs as a possible source.

SP 19-35 are located at fractures that cross McCurry Road.

Finally, SP 36-44 are located at fractures within the industrial park area.

### 4.2 HISTORICAL PHOTOGRAPHS

Table I shows the photographs that are available following a library search by National Aerial Photography for historical photographs. Since the source has not been absolutely identified, it may be prudent to review additional historical photographs that show the site from the 1950s to the 1980s. Enlargements of areas of interest may be required for future work.



## 5. DISCUSSION

The present understanding of the location of the solvent plume suggests that it is controlled by fractures. The additional sample points may further define the plume, at least to depths of 110 feet. The low levels of VOCs that have been detected make it difficult to pinpoint a source. The photographic analysis has suggested that one of the farms, present since at least 1939, could have been a possible source. Down-gradient sample locations have been recommended to explore this issue.

If businesses in the industrial park (two miles to the northeast) are indeed the responsible parties, it is very likely that the contaminants may have been transported along vertical or nearly vertical fractures, through the very permeable glacial sand and gravel deposits, and into the dolomite aquitard. In this case, VOCs may be located in pockets along the irregular dolomite surface. If this is so, then the present depth of investigation will not be sufficient to characterize the vertical extent of the plume. If it appears that VOCs, especially in the free phase, went into the groundwater, then seismic imaging will be required to define the aquitard surface and the vertical fracture conduits. This option may be discussed after the present sampling round has occurred.

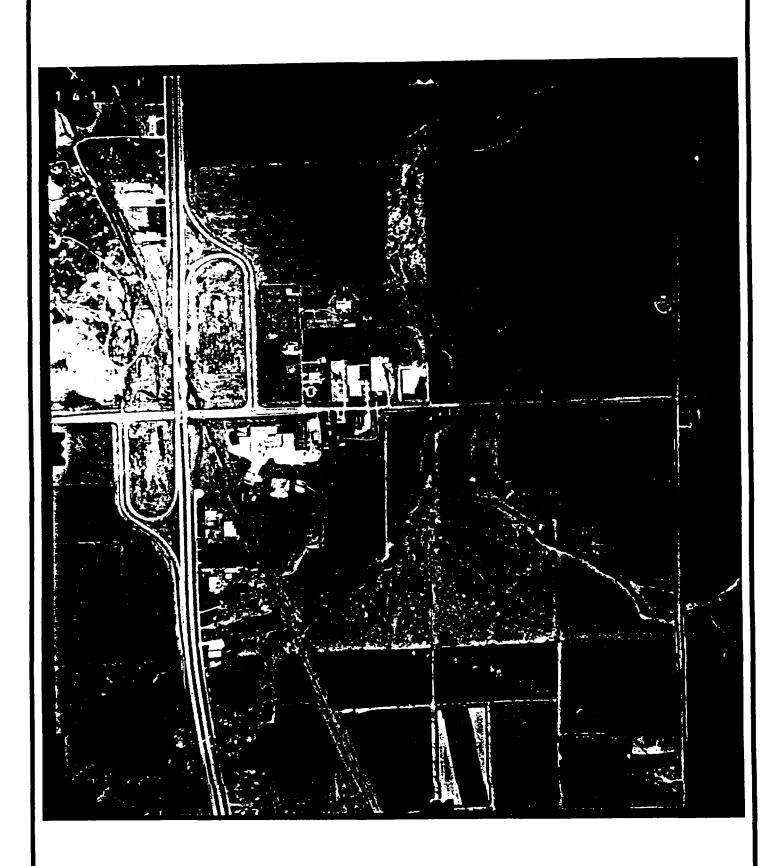
Results from the next sampling event should be reviewed and recommendations should be made, based upon this data, whether a deeper investigation is required.

## 6. REFERENCES

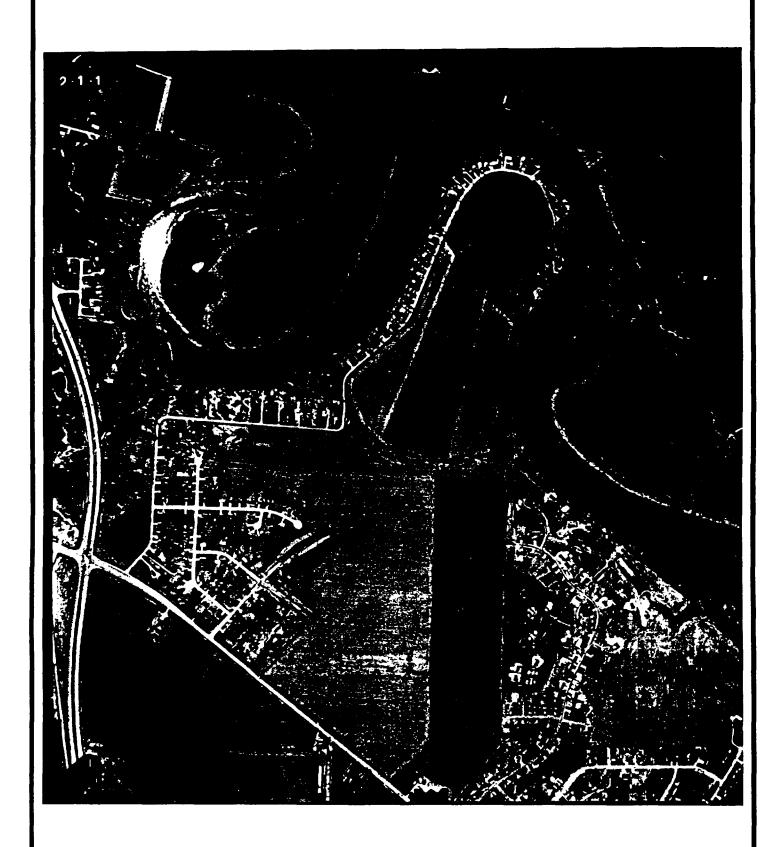
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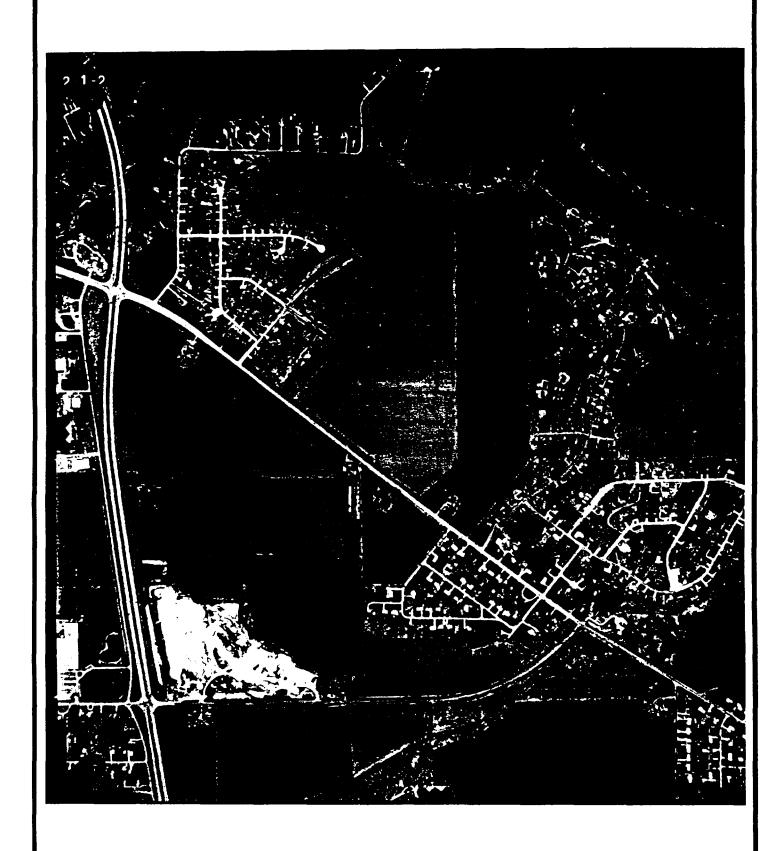
## **APPENDIX A**

UNINTERPRETED AERIAL PHOTOGRAPHS







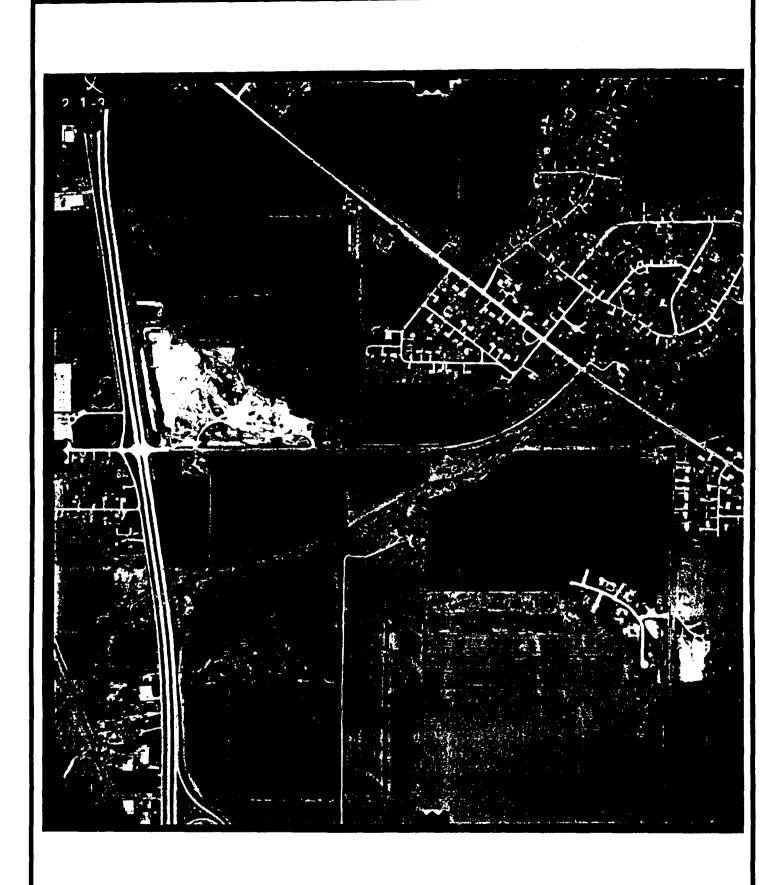


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1978 Aerial Photograph: 2-1-2

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1978 Aerial Photograph: 2-1-3

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Figure No.

